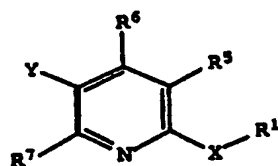




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(54) Title: SUBSTITUTED PYRIDINE COMPOUNDS AS ANTI-INFLAMMATORY AGENTS



(1)

(57) Abstract

Selected substituted pyridine compounds of formula (I) or a pharmaceutically acceptable salt thereof, wherein a. o. X is O, S, S(O), S(O)₂ or NR²; Y is -C(O)-NR³R⁴ or -NR⁴-C(O)-R³; R¹ is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of alkyl, halo, haloalkyl, cyano, azido, nitro, amidino, R¹⁸-Z¹⁸- or R¹⁸-Z¹⁸-alkyl; provided that the total number of aryl, heteroaryl, cycloalkyl and heterocyclyl radicals in R¹ is 1-3; and provided when Y is -NR⁴-C(O)-R³ and X is O or S, R¹ is other than a 2-pyrimidinyl radical; R² is a hydrogen or alkyl radical; R³ is an aryl or heteroaryl radical which is optionally substituted by 1-5 radicals of alkyl, halo, haloalkyl, cyano, azido, nitro, amidino, R¹⁹-Z¹⁹- or R¹⁹-Z¹⁹-alkyl; provided that the total number of aryl and heteroaryl radicals in R³ is 1-3; and provided when Y is -C(O)-NR³R⁴, R³ is other than a phenyl or naphthyl having an amino, nitro, cyano, carboxy or alkoxycarbonyl, are effective for prophylaxis and treatment of diseases, such as TNF-α, IL-1β, IL-6 and/or IL-8 mediated diseases, and other maladies, such as pain and diabetes. The invention encompasses novel compounds, analogs, prodrugs and pharmaceutically acceptable salts thereof, pharmaceutical compositions and methods for prophylaxis and treatment of diseases and other maladies or conditions involving inflammation, pain, diabetes, cancer and the like. The subject invention also relates to processes for making such compounds as well as to intermediates useful in such processes.

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SUBSTITUTED PYRIDINE COMPOUNDS AS ANTI-INFLAMMATORY AGENTS

BACKGROUND OF THE INVENTION

5

The present invention comprises a new class of compounds useful in treating diseases, such as TNF- α , IL-1 β , IL-6 and/or IL-8 mediated diseases and other maladies, such as pain, cancer, and diabetes. In particular, the compounds of the invention are useful for the prophylaxis and treatment of diseases or conditions involving inflammation. This invention also relates to intermediates and processes useful in the preparation of such compounds.

15 Interleukin-1 (IL-1) and Tumor Necrosis Factor α (TNF- α) are pro-inflammatory cytokines secreted by a variety of cells, including monocytes and macrophages, in response to many inflammatory stimuli (e.g., lipopolysaccharide - LPS) or external cellular stress
20 (e.g., osmotic shock and peroxide).

Elevated levels of TNF- α and/or IL-1 over basal levels have been implicated in mediating or exacerbating a number of disease states including rheumatoid arthritis; Pagets disease; osteoporosis; multiple
25 myeloma; uveitis; acute and chronic myelogenous leukemia; pancreatic β cell destruction; osteoarthritis; rheumatoid spondylitis; gouty arthritis; inflammatory bowel disease; adult respiratory distress syndrome (ARDS); psoriasis; Crohn's disease; allergic rhinitis;
30 ulcerative colitis; anaphylaxis; contact dermatitis; asthma; muscle degeneration; cachexia; Reiter's syndrome; type I and type II diabetes; bone resorption diseases; graft vs. host reaction; ischemia reperfusion injury; atherosclerosis; brain trauma; multiple
35 sclerosis; cerebral malaria; sepsis; septic shock; toxic shock syndrome; fever, and myalgias due to infection.

HIV-1, HIV-2, HIV-3, cytomegalovirus (CMV), influenza, adenovirus, the herpes viruses (including HSV-1, HSV-2), and herpes zoster are also exacerbated by TNF- α .

It has been reported that TNF- α plays a role in
5 head trauma, stroke, and ischemia. For instance, in animal models of head trauma (rat), TNF- α levels increased in the contused hemisphere (Shohami et al., *J. Cereb. Blood Flow Metab.* **14**, 615 (1994)). In a rat model of ischemia wherein the middle cerebral artery was
10 occluded, the levels of TNF- α mRNA of TNF- α increased (Feurstein et al., *Neurosci. Lett.* **164**, 125 (1993)). Administration of TNF- α into the rat cortex has been reported to result in significant neutrophil accumulation in capillaries and adherence in small blood
15 vessels. TNF- α promotes the infiltration of other cytokines (IL-1 β , IL-6) and also chemokines, which promote neutrophil infiltration into the infarct area (Feurstein, *Stroke* **25**, 1481 (1994)). TNF- α has also been implicated to play a role in type II diabetes
20 (Endocrinol. **130**, 43-52, 1994; and Endocrinol. **136**, 1474-1481, 1995).

TNF- α appears to play a role in promoting certain viral life cycles and disease states associated with them. For instance, TNF- α secreted by monocytes induced
25 elevated levels of HIV expression in a chronically infected T cell clone (Clouse et al., *J. Immunol.* **142**, 431 (1989)). Lahdevirta et al., (*Am. J. Med.* **85**, 289 (1988)) discussed the role of TNF- α in the HIV associated states of cachexia and muscle degradation.

30 TNF- α is upstream in the cytokine cascade of inflammation. As a result, elevated levels of TNF- α may lead to elevated levels of other inflammatory and proinflammatory cytokines, such as IL-1, IL-6, and IL-8.

Elevated levels of IL-1 over basal levels have been implicated in mediating or exacerbating a number of disease states including rheumatoid arthritis; osteoarthritis; rheumatoid spondylitis; gouty arthritis; inflammatory bowel disease; adult respiratory distress syndrome (ARDS); psoriasis; Crohn's disease; ulcerative colitis; anaphylaxis; muscle degeneration; cachexia; Reiter's syndrome; type I and type II diabetes; bone resorption diseases; ischemia reperfusion injury; atherosclerosis; brain trauma; multiple sclerosis; sepsis; septic shock; and toxic shock syndrome. Viruses sensitive to TNF- α inhibition, e.g., HIV-1, HIV-2, HIV-3, are also affected by IL-1.

TNF- α and IL-1 appear to play a role in pancreatic β cell destruction and diabetes. Pancreatic β cells produce insulin which helps mediate blood glucose homeostasis. Deterioration of pancreatic β cells often accompanies type I diabetes. Pancreatic β cell functional abnormalities may occur in patients with type II diabetes. Type II diabetes is characterized by a functional resistance to insulin. Further, type II diabetes is also often accompanied by elevated levels of plasma glucagon and increased rates of hepatic glucose production. Glucagon is a regulatory hormone that attenuates liver gluconeogenesis inhibition by insulin. Glucagon receptors have been found in the liver, kidney and adipose tissue. Thus glucagon antagonists are useful for attenuating plasma glucose levels (WO 97/16442, incorporated herein by reference in its entirety). By antagonizing the glucagon receptors, it is thought that insulin responsiveness in the liver will improve, thereby decreasing gluconeogenesis and lowering the rate of hepatic glucose production.

In rheumatoid arthritis models in animals, multiple intra-articular injections of IL-1 have led to an acute and destructive form of arthritis (Chandrasekhar et al., *Clinical Immunol Immunopathol.* **55**, 382 (1990)). In

studies using cultured rheumatoid synovial cells, IL-1 is a more potent inducer of stromelysin than is TNF- α (Firestein, *Am. J. Pathol.* **140**, 1309 (1992)). At sites of local injection, neutrophil, lymphocyte, and monocyte emigration has been observed. The emigration is attributed to the induction of chemokines (e.g., IL-8), and the up-regulation of adhesion molecules (Dinarello, *Eur. Cytokine Netw.* **5**, 517-531 (1994)).

IL-1 also appears to play a role in promoting certain viral life cycles. For example, cytokine-induced increase of HIV expression in a chronically infected macrophage line has been associated with a concomitant and selective increase in IL-1 production (Folks et al., *J. Immunol.* **136**, 40 (1986)). Beutler et al. (*J. Immunol.* **135**, 3969 (1985)) discussed the role of IL-1 in cachexia. Baracos et al. (*New Eng. J. Med.* **308**, 553 (1983)) discussed the role of IL-1 in muscle degeneration.

In rheumatoid arthritis, both IL-1 and TNF- α induce synoviocytes and chondrocytes to produce collagenase and neutral proteases, which leads to tissue destruction within the arthritic joints. In a model of arthritis (collagen-induced arthritis (CIA) in rats and mice), intra-articular administration of TNF- α either prior to or after the induction of CIA led to an accelerated onset of arthritis and a more severe course of the disease (Brahm et al., *Lymphokine Cytokine Res.* **11**, 253 (1992); and Cooper, *Clin. Exp. Immunol.* **898**, 244 (1992)).

IL-8 has been implicated in exacerbating and/or causing many disease states in which massive neutrophil infiltration into sites of inflammation or injury (e.g., ischemia) is mediated by the chemotactic nature of IL-8, including, but not limited to, the following: asthma, inflammatory bowel disease, psoriasis, adult respiratory distress syndrome, cardiac and renal reperfusion injury,

thrombosis and glomerulonephritis. In addition to the chemotaxis effect on neutrophils, IL-8 also has the ability to activate neutrophils. Thus, reduction in IL-8 levels may lead to diminished neutrophil infiltration.

5 Several approaches have been taken to block the effect of TNF- α . One approach involves using soluble receptors for TNF- α (e.g., TNFR-55 or TNFR-75), which have demonstrated efficacy in animal models of TNF- α -mediated disease states. A second approach to
10 neutralizing TNF- α using a monoclonal antibody specific to TNF- α , cA2, has demonstrated improvement in swollen joint count in a Phase II human trial of rheumatoid arthritis (Feldmann et al., *Immunological Reviews*, pp. 195-223 (1995)). These approaches block the effects of
15 TNF- α and IL-1 by either protein sequestration or receptor antagonism.

 The present invention also relates to a method of treating cancer which is mediated by Raf and Raf-inducible proteins. Raf proteins are kinases activated
20 in response to extracellular mitogenic stimuli such as PDGF, EGF, acidic FGF, thrombin, insulin or endothelin, and also in response to oncoproteins such as v-src, v-sis, and v-fms. Raf functions downstream of ras in signal transduction from the cellular membrane to the
25 nucleus. Compounds in the present invention may be oncolytics through the antagonism of Raf kinase. Antisense constructs which reduce cellular levels of c-Raf and hence Raf activity inhibit the growth of rodent fibroblasts in soft agar, while exhibiting little or no
30 general cytotoxicity. This inhibition of growth in soft agar is highly predictive of tumor responsiveness in whole animals. Moreover Raf antisense constructs have shown efficacy in reducing tumor burden in animals. Examples of cancers where Raf kinase is implicated by
35 overexpression include cancers of the brain, larynx, lung, lymphatic system, urinary tract and stomach,

including hystocytic lymphoma, lung adenocarcinoma and small cell lung cancers. Other examples include cancers involving overexpression of upstream activators of Raf or Raf-activating oncogenes, including pancreatic and
5 breast carcinoma.

Substituted imidazole and pyrrole compounds have been described for use in the treatment of cytokine mediated diseases by inhibition of proinflammatory cytokines, such as IL-1, IL-6, IL-8 and TNF.
10 Substituted imidazoles for use in the treatment of cytokine mediated diseases have been described in US Pat. 5,593,992; WO 93/14081; WO 97/18626; WO 96/21452; WO 96/21654; WO 96/40143; WO 97/05878; WO 97/05878; (each of which is incorporated herein by reference in
15 its entirety). Substituted imidazoles for use in the treatment of inflammation has been described in US Pat. 3,929,807 (which is incorporated herein by reference in its entirety). Substituted pyrrole compounds for use in the treatment of cytokine mediated diseases have been
20 described in WO 97/05877; WO 97/05878; WO 97/16426; WO 97/16441; and WO 97/16442 (each of which is incorporated herein by reference in its entirety).

Substituted 2-aminopyridine compounds have been described as nitric oxide synthase inhibitors for use in
25 the treatment of inflammation, neurodegenerative disorders and disorders of gastrointestinal motility in WO 96/18616 and WO 96/18617.

Diaryl substituted pyridine compounds have been described for use in the treatment of inflammation and
30 inflammation related disorders in WO 96/24584 and US 5,596,008.

US 3,980,652, US 3,991,057 and US 4,002,629 describe piperazinyl substituted pyridine compounds for use as anti-inflammatory and cardiovascular agents.

35 JP 6135934 describes substituted pyridine compounds as phospholipase A2 inhibitors for use as antiphlogistic and anti-pancreatitis agents. GB 1,189,188 describes

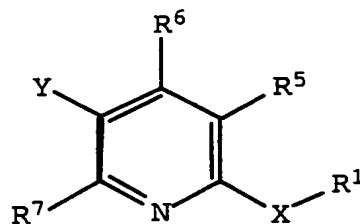
pyrimidin-2-ylamino substituted pyridine compounds as therapeutically valuable compounds for use as antiphlogistic agents.

5

BRIEF DESCRIPTION OF THE INVENTION

The present invention comprises a new class of compounds useful in the prophylaxis and treatment of diseases, such as TNF- α , IL-1 β , IL-6 and/or IL-8 mediated diseases and other maladies, such as pain, cancer, and diabetes. In particular, the compounds of the invention are useful for the prophylaxis and treatment of diseases or conditions involving inflammation. Accordingly, the invention also comprises pharmaceutical compositions comprising the compounds, methods for the prophylaxis and treatment of TNF- α , IL-1 β , IL-6 and/or IL-8 mediated diseases, such as inflammatory, pain and diabetes diseases, using the compounds and compositions of the invention, and intermediates and processes useful for the preparation of the compounds of the invention.

The compounds of the invention are represented by the following general structure:



25

(I)

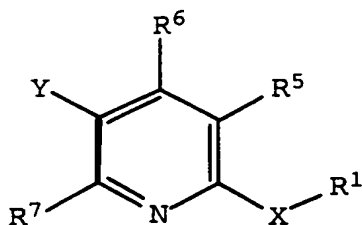
wherein R¹, R⁵, R⁶, R⁷, X and Y are defined below.

The foregoing merely summarizes certain aspects of the invention and is not intended, nor should it be construed, as limiting the invention in any way. All patents and other publications recited herein are hereby incorporated by reference in their entirety.

30

DETAILED DESCRIPTION OF THE INVENTION

In accordance with the present invention, there is provided compounds of the formula:



(I)

or a pharmaceutically acceptable salt thereof, wherein

X is O, S, S(O), S(O)₂, or NR²; preferably, X is O, S or NR²; more preferably, X is O or NR²; most preferably, X is NR²;

Y is -C(O)-NR³R⁴ or -NR⁴-C(O)-R³;

R¹ is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of alkyl, halo, haloalkyl, cyano, azido, nitro, amidino, R¹⁸-Z¹⁸- or R¹⁸-Z¹⁸-alkyl;

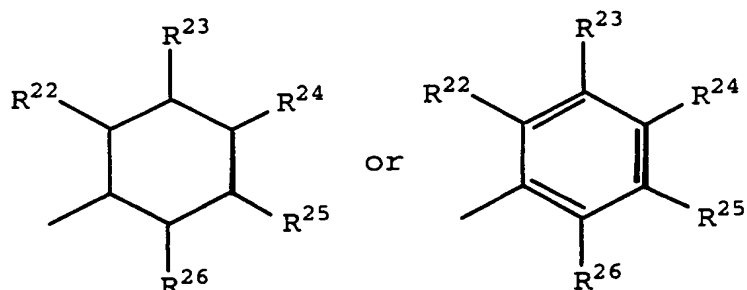
preferably, R¹ is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of C₁-C₆ alkyl, halo, C₁-C₆ haloalkyl of 1-3 halo radicals, cyano, azido, nitro, amidino, R¹⁸-Z¹⁸- or R¹⁸-Z¹⁸-C₁-C₆ alkyl;

more preferably, R¹ is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of C₁-C₄ alkyl, halo, C₁-C₄ haloalkyl of 1-3

halo radicals, cyano, azido, nitro, amidino, $R^{18}-Z^{18}-$ or $R^{18}-Z^{18}-C_1-C_4$ alkyl;

provided that the total number of aryl, heteroaryl,
 5 cycloalkyl and heterocyclyl radicals in R^1 is 1-3,
 preferably, 1-2, and provided when Y is $-NR^4-C(O)-R^3$ and
 X is O or S, R^1 is other than a 2-pyrimidinyl radical;

more preferably, R^1 is a radical of the formula



10

wherein R^{22} , R^{23} , R^{24} , R^{25} and R^{26} are each independently a
 radical of hydrogen, C_1-C_4 alkyl, halo, trifluoromethyl,
 cyano, azido, nitro, amidino, $R^{18}-Z^{18}-$ or $R^{18}-Z^{18}-C_1-C_4$
 alkyl; provided at least one of R^{21} , R^{22} , R^{23} , R^{24} and R^{25}
 15 is hydrogen; and provided that the combined total number
 of aryl and heteroaryl radicals in R^{22} , R^{23} , R^{24} , R^{25} and
 R^{26} is 0-1;

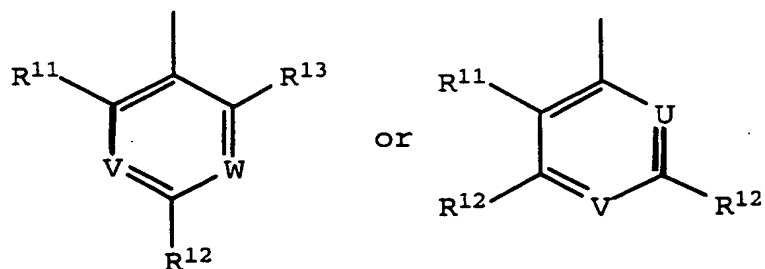
R^2 is a hydrogen or alkyl radical; preferably, R^2 is a
 20 hydrogen or C_1-C_4 alkyl radical; more preferably, R^2 is a
 hydrogen or C_1-C_2 alkyl radical; more preferably, R^2 is a
 hydrogen or methyl radical; and most preferably, R^2 is a
 hydrogen radical;

25 R^3 is an aryl or heteroaryl radical which is optionally
 substituted by 1-5 radicals of alkyl, halo, haloalkyl,

cyano, azido, nitro, amidino, $R^{19}-Z^{19}$ - or $R^{19}-Z^{19}$ -alkyl;
 preferably, R^3 is an aryl or heteroaryl radical which is
 optionally substituted by 1-5 radicals of C_1-C_6 alkyl,
 halo, C_1-C_6 haloalkyl of 1-3 halo radicals, cyano,
 5 azido, nitro, amidino, $R^{19}-Z^{19}$ - or $R^{19}-Z^{19}-C_1-C_6$ alkyl;
 more preferably, R^3 is an aryl or heteroaryl radical
 which is optionally substituted by 1-5 radicals of C_1-C_6
 alkyl, halo, C_1-C_4 haloalkyl of 1-3 halo radicals,
 cyano, azido, nitro, amidino, $R^{19}-Z^{19}$ - or $R^{19}-Z^{19}-C_1-C_4$
 10 alkyl;

provided that the total number of aryl and heteroaryl
 radicals in R^3 is 1-3, preferably, 1-2; and provided
 when Y is $-C(O)-NR^3R^4$, R^3 is other than a phenyl or
 15 naphthyl having an amino, nitro, cyano, carboxy or
 alkoxy carbonyl substituent bonded to the ring carbon
 atom adjacent to the ring carbon atom bonded to $-NR^4$;

more preferably, R^3 is a radical of the formula



20

wherein

U is $C-R^{13}$ or N;

25 V and W are each independently $C-R^{12}$ or N;

R^{11} and R^{13} are each independently a radical of hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl, cyano, azido, nitro, amidino or $R^{19}-Z^{19}-$; preferably, R^{11} and R^{13} are each independently a radical of hydrogen, methyl, ethyl, fluoro, chloro, trifluoromethyl, cyano, azido, nitro, amidino, $R^{19}-O-$, $R^{19}-S(O)_2-$, $R^{19}-O-C(O)-$, $R^{19}-C(O)-$, $R^{19}-NR^{21}-C(O)-$ or $R^{19}-NR^{21}-S(O)_2-$;

each R^{12} is independently a radical of hydrogen, C₁-C₆ alkyl, halo, C₁-C₄ haloalkyl of 1-3 halo radicals, $R^{31}-Z^{31}-$ or $R^{31}-Z^{31}-C_1-C_4$ alkyl; preferably, each R^{12} is independently a radical of hydrogen, methyl, ethyl, fluoro, chloro, trifluoromethyl, trifluoromethoxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, aminocarbonyl, methylaminocarbonyl, dimethylaminocarbonyl, aminomethyl, (methylamino)methyl or (dimethylamino)methyl;

provided that the combined total number of aryl and heteroaryl radicals in R^{11} , R^{12} and R^{13} is 0-1;

wherein each R^{31} is independently a hydrogen, C₁-C₄ alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy;

each Z^{31} is independently $-O-$, $-NR^{21}-$, $-NR^{21}-C(O)-$, $-C(O)-NR^{21}-$, $-NR^{21}-S(O)_2-$ or $-S(O)_2-NR^{21}-$;

- R^4 is a hydrogen, alkyl, alkenyl, haloalkyl, haloalkenyl, aryl, heteroaryl, arylalkyl, heteroarylalkyl or $R^{20}-Z^{20}$ -alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, alkoxy, alkylthiol, amino, alkylamino, dialkylamino, alkanoylamino, alkylsulfonylamino, alkylsulfinyl, alkylsulfonyl, alkoxycarbonylamino, alkoxycarbonyl, cyano, halo, azido, alkyl, haloalkyl or haloalkoxy;
- preferably, R^4 is a radical of hydrogen, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_1 - C_6 haloalkyl of 1-3 halo radicals, C_2 - C_6 haloalkenyl of 1-3 halo radicals, aryl, heteroaryl, aryl- C_1 - C_4 alkyl, heteroaryl- C_1 - C_4 alkyl or $R^{20}-Z^{20}$ - C_1 - C_6 alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, C_1 - C_5 alkanoylamino, C_1 - C_4 alkylsulfonylamino, C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkylsulfonyl, (C_1 - C_4 alkoxy)carbonylamino, (C_1 - C_4 alkoxy)carbonyl, cyano, halo, azido, C_1 - C_4 alkyl, C_1 - C_4 haloalkyl of 1-3 halo radicals or C_1 - C_4 haloalkoxy of 1-3 halo radicals;
- more preferably, R^4 is a radical of hydrogen, C_1 - C_6 alkyl, aryl, heteroaryl, aryl- C_1 - C_4 alkyl, heteroaryl- C_1 - C_4 alkyl or $R^{20}-Z^{20}$ - C_2 - C_4 alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, acetylamino, halo, C_1 - C_4 alkyl, trifluoromethyl or trifluoromethoxy;

more preferably, R^4 is a radical of hydrogen, C_1 - C_6 alkyl, aryl, heteroaryl, aryl- C_1 - C_4 alkyl, heteroaryl- C_1 - C_4 alkyl or R^{20} - Z^{20} - C_2 - C_4 alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted
5 by 1-2 radicals of hydroxy, methoxy, ethoxy, methylthiol, ethylthiol, amino, methylamino, dimethylamino, ethylamino, diethylamino, acetylamino, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy;

10

more preferably, R^4 is a radical of hydrogen, methyl or ethyl radical;

wherein each R^{18} is independently a hydrogen, alkyl,
15 haloalkyl, aryl, heteroaryl, arylalkyl or heteroarylalkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, alkoxy, alkylthiol, amino, alkylamino, dialkylamino, alkanoylamino, alkylsulfonylamino,
20 alkylsulfinyl, alkylsulfonyl, alkoxycarbonylamino, alkoxycarbonyl, cyano, halo, azido, alkyl, haloalkyl or haloalkoxy;

preferably, each R^{18} is independently a hydrogen, C_1 - C_4
25 alkyl, C_1 - C_4 haloalkyl of 1-3 halo radicals, aryl, heteroaryl, aryl- C_1 - C_4 alkyl or heteroaryl- C_1 - C_4 alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 -
30 C_4 alkyl)amino, C_1 - C_5 alkanoylamino, C_1 - C_4 alkylsulfonylamino, C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkylsulfonyl, (C_1 - C_4 alkoxy)carbonylamino, (C_1 - C_4 alkoxy)carbonyl, cyano, halo, azido, C_1 - C_4 alkyl, C_1 - C_4 haloalkyl of 1-3 halo radicals or C_1 - C_4 haloalkoxy of 1-
35 3 halo radicals;

more preferably, each R^{18} is independently a hydrogen, C_1 - C_4 alkyl, trifluoromethyl, aryl, heteroaryl, aryl- C_1 - C_2 alkyl or heteroaryl- C_1 - C_2 alkyl radical, wherein the
5 aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, acetylamino, cyano, halo, azido, C_1 - C_4 alkyl, trifluoromethyl or trifluoromethoxy;
10 each Z^{18} is independently -O-, -S-, -S(O)-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; preferably, each Z^{18} is independently -O-, -S-, -S(O)₂-, -CO₂-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-;
15

wherein each R^{19} is independently a hydrogen, alkyl, haloalkyl, aryl, heteroaryl, arylalkyl or heteroarylalkyl radical, wherein the aryl and heteroaryl
20 radicals are optionally substituted by 1-3 radicals of hydroxy, alkoxy, alkylthiol, amino, alkylamino, dialkylamino, alkanoylamino, alkylsulfonylamino, alkylsulfinyl, alkylsulfonyl, alkoxycarbonylamino, alkoxycarbonyl, cyano, halo, azido, alkyl, haloalkyl or
25 haloalkoxy;

preferably, each R^{19} is independently a hydrogen, C_1 - C_4 alkyl, C_1 - C_4 haloalkyl of 1-3 halo radicals, aryl, heteroaryl, aryl- C_1 - C_4 alkyl or heteroaryl- C_1 - C_4 alkyl
30 radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, C_1 - C_5 alkanoylamino, C_1 - C_4 alkylsulfonylamino, C_1 - C_4 alkylsulfinyl, C_1 - C_4

alkylsulfonyl, (C₁-C₄ alkoxy)carbonylamino, (C₁-C₄ alkoxy)carbonyl, cyano, halo, azido, C₁-C₄ alkyl, C₁-C₄ haloalkyl of 1-3 halo radicals or C₁-C₄ haloalkoxy of 1-3 halo radicals;

5

more preferably, each R¹⁹ is independently a hydrogen, C₁-C₄ alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted
 10 by 1-2 radicals of hydroxy, C₁-C₄ alkoxy, C₁-C₄ alkylthiol, amino, C₁-C₄ alkylamino, di(C₁-C₄ alkyl)amino, acetylamino, cyano, halo, C₁-C₄ alkyl, trifluoromethyl or trifluoromethoxy;

15 more preferably, each R¹⁹ is independently a hydrogen, C₁-C₄ alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino,
 20 methylamino, dimethylamino, acetylamino, cyano, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy;

more preferably, each R¹⁹ is independently a hydrogen, methyl, ethyl, trifluoromethyl, phenyl, heteroaryl,
 25 phenylmethyl or heteroaryl-methyl radical, wherein the phenyl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, fluoro, chloro, methyl, ethyl, trifluoromethyl or
 30 trifluoromethoxy;

each Z¹⁹ is independently -O-, -S-, -S(O)-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; preferably, each Z¹⁹ is

independently -O-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; more preferably, each Z¹⁹ is independently -O-, -S(O)₂-, -O-C(O)-, -C(O)-, -NR²¹-C(O)- or -NR²¹-S(O)₂-;

5

wherein each R²⁰ is independently a hydrogen, alkyl, haloalkyl, aryl, heteroaryl, arylalkyl or heteroarylalkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, alkoxy, alkylthiol, amino, alkylamino, dialkylamino, alkanoylamino, alkylsulfonylamino, alkylsulfinyl, alkylsulfonyl, alkoxycarbonylamino, alkoxycarbonyl, cyano, halo, azido, alkyl, haloalkyl or haloalkoxy;

15

preferably, each R²⁰ is independently a hydrogen, C₁-C₄ alkyl, C₁-C₄ haloalkyl of 1-3 halo radicals, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-3 radicals of hydroxy, C₁-C₄ alkoxy, C₁-C₄ alkylthiol, amino, C₁-C₄ alkylamino, di(C₁-C₄ alkyl)amino, C₁-C₅ alkanoylamino, C₁-C₄ alkylsulfonylamino, C₁-C₄ alkylsulfinyl, C₁-C₄ alkylsulfonyl, (C₁-C₄ alkoxy)carbonylamino, (C₁-C₄ alkoxy)carbonyl, cyano, halo, azido, C₁-C₄ alkyl, C₁-C₄ haloalkyl of 1-3 halo radicals or C₁-C₄ haloalkoxy of 1-3 halo radicals;

more preferably, each R²⁰ is independently a hydrogen, C₁-C₄ alkyl, aryl, heteroaryl, aryl-C₁-C₂ alkyl or heteroaryl-C₁-C₂ alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, C₁-C₄ alkoxy, C₁-C₄ alkylthiol, amino, C₁-C₄ alkylamino, di(C₁-C₄ alkyl)amino,

30

acetylamino, halo, C₁-C₄ alkyl, trifluoromethyl or trifluoromethoxy;

more preferably, each R²⁰ is independently a hydrogen,
5 C₁-C₄ alkyl, aryl, heteroaryl, aryl-C₁-C₂ alkyl or heteroaryl-C₁-C₂ alkyl radical, wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, methylthiol, ethylthiol, amino, methylamino, dimethylamino,
10 ethylamino, diethylamino, acetylamino, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy;

each Z²⁰ is independently -O-, -S-, -S(O)-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-
15 S(O)₂- or -S(O)₂-NR²¹-; preferably, each Z²⁰ is independently -O- or -NR²¹-;

wherein each R²¹ is independently a hydrogen or alkyl radical; preferably, each R²¹ is independently a hydrogen
20 or C₁-C₄ alkyl radical; more preferably, each R²¹ is independently a hydrogen or methyl radical;

R⁵ and R⁶ are each independently a hydrogen, alkyl, halo, haloalkyl, haloalkoxy, aminoalkyl, alkylaminoalkyl,
25 dialkylaminoalkyl, amino, alkylamino, dialkylamino, alkanoylamino, alkylsulfonylamino, aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, hydroxy, hydroxyalkyl, thiol, alkylthiol, alkylsulfinyl, alkylsulfonyl, alkoxy, alkoxyalkyl, cyano, azido, nitro,
30 carboxy, alkoxycarbonyl, aminocarbonyl, alkylaminocarbonyl or dialkylaminocarbonyl radical;

preferably, R⁵ and R⁶ are each independently a hydrogen, C₁-C₄ alkyl, halo, C₁-C₄ haloalkyl of 1-3 halo radicals,

C₁-C₄ haloalkoxy of 1-3 halo radicals, C₁-C₄ aminoalkyl, (C₁-C₄ alkyl)amino-C₁-C₄ alkyl, di(C₁-C₄ alkyl)amino-C₁-C₄ alkyl, amino, C₁-C₄ alkylamino, di(C₁-C₄ alkyl)amino, C₁-C₅ alkanoylamino, C₁-C₄ alkylsulfonylamino, 5 aminosulfonyl, C₁-C₄ alkylaminosulfonyl, di(C₁-C₄ alkyl)aminosulfonyl, hydroxy, C₁-C₄ hydroxyalkyl, thiol, C₁-C₄ alkylthiol, C₁-C₄ alkylsulfinyl, C₁-C₄ alkylsulfonyl, C₁-C₄ alkoxy, (C₁-C₄ alkoxy)C₁-C₄ alkyl, cyano, azido, nitro, carboxy, (C₁-C₄ alkoxy)carbonyl, 10 aminocarbonyl, (C₁-C₄ alkyl)aminocarbonyl or di(C₁-C₄ alkyl)aminocarbonyl radical;

more preferably, R⁵ and R⁶ are each independently a hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl, 15 trifluoromethoxy, amino, C₁-C₄ alkylamino, di(C₁-C₄ alkyl)amino, C₁-C₅ alkanoylamino, hydroxy, C₁-C₄ hydroxyalkyl, C₁-C₄ alkoxy, cyano, azido, nitro, carboxy, (C₁-C₄ alkoxy)carbonyl, aminocarbonyl, (C₁-C₄ alkyl)aminocarbonyl or di(C₁-C₄ alkyl)aminocarbonyl 20 radical;

more preferably, R⁵ and R⁶ are each independently a hydrogen, methyl, ethyl, halo, trifluoromethyl, trifluoromethoxy, amino, C₁-C₂ alkylamino, di(C₁-C₂ 25 alkyl)amino, hydroxy, methoxy or ethoxy radical; most preferably, R⁵ and R⁶ are each a hydrogen radical;

R⁷ is a hydrogen, alkyl, halo, haloalkyl, haloalkoxy, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, 30 aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, hydroxy, hydroxyalkyl, thiol, alkylthiol, alkylsulfinyl, alkylsulfonyl, alkoxy, alkoxyalkyl, cyano, azido, nitro, carboxy, alkoxycarbonyl, aminocarbonyl, alkylaminocarbonyl or dialkylaminocarbonyl radical;

preferably, R^7 is a hydrogen, C_1 - C_4 alkyl, halo, C_1 - C_4 haloalkyl of 1-3 halo radicals, C_1 - C_4 haloalkoxy of 1-3 halo radicals, C_1 - C_4 aminoalkyl, (C_1 - C_4 alkyl)amino- C_1 - C_4 alkyl, di(C_1 - C_4 alkyl)amino- C_1 - C_4 alkyl, aminosulfonyl, C_1 - C_4 alkylaminosulfonyl, di(C_1 - C_4 alkyl)aminosulfonyl, hydroxy, C_1 - C_4 hydroxyalkyl, thiol, C_1 - C_4 alkylthiol, C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkylsulfonyl, C_1 - C_4 alkoxy, (C_1 - C_4 alkoxy) C_1 - C_4 alkyl, cyano, azido, nitro, carboxy, (C_1 - C_4 alkoxy)carbonyl, aminocarbonyl, (C_1 - C_4 alkyl)aminocarbonyl or di(C_1 - C_4 alkyl)aminocarbonyl radical;

more preferably, R^7 is a hydrogen, C_1 - C_4 alkyl, halo, trifluoromethyl, trifluoromethoxy, hydroxy, C_1 - C_4 hydroxyalkyl, C_1 - C_4 alkoxy, carboxy, (C_1 - C_4 alkoxy)carbonyl, aminocarbonyl, (C_1 - C_4 alkyl)aminocarbonyl or di(C_1 - C_4 alkyl)aminocarbonyl radical;

more preferably, R^7 is a hydrogen, methyl, ethyl, halo, trifluoromethyl, trifluoromethoxy, hydroxy, methoxy or ethoxy radical; most preferably, R^7 is a hydrogen radical.

The compounds of this invention may have in general several asymmetric centers and are typically depicted in the form of racemic mixtures. This invention is intended to encompass racemic mixtures, partially racemic mixtures and separate enantiomers and diastereomers.

Compounds of interest include the following:
2-cyclohexyloxy-5-(2-chlorophenylcarbonylamino)pyridine;
2-cyclohexyloxy-5-(2-methylphenylcarbonylamino)pyridine;

- 2-cyclohexyloxy-5-(2,6-dichlorophenylcarbonylamino)
pyridine;
2-cyclohexyloxy-5-(2,6-dimethylphenylcarbonylamino)
pyridine;
- 5 2-(2,4-dimethylphenoxy)-5-(2-chlorophenylcarbonylamino)
pyridine;
2-(2,4-dimethylphenoxy)-5-(2,6-dichlorophenylcarbonyl
amino)pyridine;
2-(2,4-dimethylphenoxy)-5-(2-methylphenylcarbonylamino)
- 10 pyridine;
2-(2,6-dimethyl-4-chlorophenoxy)-5-(2,6-dimethylphenyl
carbonylamino) pyridine;
2-(2-methyl-4-fluorophenoxy)-5-(2-methylphenylcarbonyl
amino)pyridine;
- 15 2-(2-methyl-4-chlorophenoxy)-5-(2-chlorophenylcarbonyl
amino)pyridine;
2-(2-methyl-4-chlorophenoxy)-5-(2-methylphenylcarbonyl
amino)pyridine;
2-(2-methylphenoxy)-5-(2-chlorophenylcarbonylamino)
- 20 pyridine;
2-(2-methylphenoxy)-5-(2,6-dichlorophenyl
carbonylamino)pyridine;
2-(2-methylphenoxy)-5-(2-methylphenylcarbonyl
amino)pyridine;
- 25 2-(2-methyl-4-chlorophenoxy)-5-(2,6-dichlorophenyl
carbonylamino)pyridine;
2-(2-methyl-4-chlorophenoxy)-5-(2,6-dimethylphenyl
carbonylamino)pyridine;
2-(4-chlorophenoxy)-5-(2,6-dimethylphenylcarbonylamino)
- 30 pyridine;
2-(2-methyl-4-fluorophenoxy)-5-(2,6-dichlorophenyl
carbonylamino)pyridine;
2-(2-methyl-4-fluorophenoxy)-5-(2,6-dimethylphenyl
carbonylamino)pyridine;
- 35 2-(2-methylphenoxy)-5-(2,6-dimethylphenyl
carbonylamino)pyridine;

- 2-(2-methyl-4-fluorophenoxy)-5-(2-fluorophenylcarbonyl amino)pyridine;
2-(2,4-dimethylphenoxy)-5-(2,6-dimethylphenylcarbonyl amino)pyridine;
5 2-(1-naphthyloxy)-5-(2-methylphenylcarbonylamino) pyridine;
2-(1-naphthyloxy)-5-(2,6-dichlorophenylcarbonylamino) pyridine;
2-(1-naphthyloxy)-5-(2,6-dimethylphenylcarbonylamino) 10 pyridine;
2-(2-methyl-3-pyridyloxy)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
2-(2-methyl-4-chlorophenoxy)-5-((3,5-dimethyl-4-isoxazolyl)carbonylamino)pyridine;
15 2-(2-methyl-4-chlorophenylthiol)-5-(2-methylphenylcarbonyl amino)pyridine;
2-(2-methyl-4-chlorophenylthiol)-5-(2,6-dimethylphenylcarbonyl amino)pyridine;
2-cyclohexylamino-5-(2,6-dichlorophenylcarbonylamino) 20 pyridine;
2-cyclohexylamino-5-(2,6-dimethylphenylcarbonylamino) pyridine;
2-(2-methylcyclohexylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
25 2-(2-methylcyclohexylamino)-5-(2-methylphenylcarbonyl amino)pyridine;
2-(2,4-dimethylphenylamino)-5-(2-fluorophenylcarbonyl amino)pyridine;
2-(2,4-dimethylphenylamino)-5-(2-chlorophenylcarbonyl 30 amino)pyridine;
2-(2,4-dimethylphenylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
2-(2-methyl-4-chlorophenylamino)-5-(2,6-dichlorophenylcarbonylamino)pyridine;
35 2-(2,4-dimethylphenylamino)-5-(2-methylphenylcarbonyl amino)pyridine;

2-(2-methylphenylamino)-5-(2-methylphenylcarbonyl
amino)pyridine;
2-(2-methylphenylamino)-5-(2,6-dichlorophenylcarbonyl
amino)pyridine;
5 2-(2-methylphenylamino)-5-(2,6-dimethylphenylcarbonyl
amino)pyridine;
2-(2,4-dimethylphenylamino)-5-(2,6-
dimethylphenylcarbonyl amino)pyridine;
2-(2-methyl-4-chlorophenylamino)-5-(2-methylphenyl
10 carbonylamino)pyridine;
2-(2-methyl-4-chlorophenylamino)-5-(2,6-dimethylphenyl
carbonylamino)pyridine; and
2-(2-methyl-4-chlorophenylamino)-5-(2-methylphenyl
aminocarbonyl)pyridine.

15

As utilized herein, the following terms shall have
the following meanings:

"Alkyl", alone or in combination, means a straight-chain
20 or branched-chain alkyl radical containing preferably 1-
15 carbon atoms (C_1 - C_{15}), more preferably 1-8 carbon
atoms (C_1 - C_8), even more preferably 1-6 carbon atoms
(C_1 - C_6), yet more preferably 1-4 carbon atoms (C_1 - C_4),
still more preferably 1-3 carbon atoms (C_1 - C_3), and most
25 preferably 1-2 carbon atoms (C_1 - C_2). Examples of such
radicals include methyl, ethyl, n-propyl, isopropyl,
n-butyl, isobutyl, sec-butyl, tert-butyl, pentyl, iso-
amyl, hexyl, octyl and the like.

30 "Hydroxyalkyl", alone or in combination, means an alkyl
radical as defined above wherein at least one hydrogen
radical is replaced with a hydroxyl radical, preferably
1-3 hydrogen radicals are replaced by hydroxyl radicals,
more preferably 1-2 hydrogen radicals are replaced by
35 hydroxyl radicals, and most preferably one hydrogen
radical is replaced by a hydroxyl radical. Examples of
such radicals include hydroxymethyl, 1-, 2-hydroxyethyl,

1-, 2-, 3-hydroxypropyl, 1,3-dihydroxy-2-propyl, 1,3-dihydroxybutyl, 1,2,3,4,5,6-hexahydroxy-2-hexyl and the like.

5 "Alkenyl", alone or in combination, means a straight-chain or branched-chain hydrocarbon radical having one or more double bonds, preferably 1-2 double bonds and more preferably one double bond, and containing preferably 2-15 carbon atoms (C_2-C_{15}), more preferably
10 2-8 carbon atoms (C_2-C_8), even more preferably 2-6 carbon atoms (C_2-C_6), yet more preferably 2-4 carbon atoms (C_2-C_4), and still more preferably 2-3 carbon atoms (C_2-C_3). Examples of such alkenyl radicals include ethenyl, propenyl, 2-methylpropenyl, 1,4-
15 butadienyl and the like.

"Alkoxy", alone or in combination, means a radical of the type "R-O-" wherein "R" is an alkyl radical as defined above and "O" is an oxygen atom. Examples of
20 such alkoxy radicals include methoxy, ethoxy, n-propoxy, isopropoxy, n-butoxy, iso-butoxy, sec-butoxy, tert-butoxy and the like.

"Alkoxycarbonyl", alone or in combination, means a
25 radical of the type "R-O-C(O)-" wherein "R-O-" is an alkoxy radical as defined above and "C(O)" is a carbonyl radical.

"Alkoxycarbonylamino", alone or in combination, means a
30 radical of the type "R-O-C(O)-NH-" wherein "R-O-C(O)" is an alkoxycarbonyl radical as defined above, wherein the amino radical may optionally be substituted, such as with alkyl, aryl, aralkyl, cycloalkyl, cycloalkylalkyl and the like.

35 "Alkylthio", alone or in combination, means a radical of the type "R-S-" wherein "R" is an alkyl radical as

defined above and "S" is a sulfur atom. Examples of such alkylthio radicals include methylthio, ethylthio, n-propylthio, isopropylthio, n-butylthio, iso-butylthio, sec-butylthio, tert-butylthio and the like.

5

"Alkylsulfinyl", alone or in combination, means a radical of the type "R-S(O)-" wherein "R" is an alkyl radical as defined above and "S(O)" is a mono-oxygenated sulfur atom. Examples of such alkylsulfinyl radicals include methylsulfinyl, ethylsulfinyl, n-propylsulfinyl, isopropylsulfinyl, n-butylsulfinyl, iso-butylsulfinyl, sec-butylsulfinyl, tert-butylsulfinyl and the like.

"Alkylsulfonyl", alone or in combination, means a radical of the type "R-S(O)₂-" wherein "R" is an alkyl radical as defined above and "S(O)₂" is a di-oxygenated sulfur atom. Examples of such alkylsulfonyl radicals include methylsulfonyl, ethylsulfonyl, n-propylsulfonyl, isopropylsulfonyl, n-butylsulfonyl, iso-butylsulfonyl, sec-butylsulfonyl, tert-butylsulfonyl and the like.

20

"Aryl", alone or in combination, means a phenyl or biphenyl radical, which is optionally benzo fused or heterocyclo fused and which is optionally substituted with one or more substituents selected from alkyl, alkoxy, halogen, hydroxy, amino, azido, nitro, cyano, haloalkyl, carboxy, alkoxycarbonyl, cycloalkyl, alkanoylamino, amido, amidino, alkoxycarbonylamino, N-alkylamidino, alkylamino, dialkylamino, aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl, N-alkylamido, N,N-dialkylamido, aralkoxycarbonylamino, alkylthio, alkylsulfinyl, alkylsulfonyl, oxo and the like. Examples of aryl radicals are phenyl, o-tolyl, 4-methoxyphenyl, 2-(tert-butoxy)phenyl, 3-methyl-4-methoxyphenyl, 2-CF₃-phenyl, 2-fluorophenyl, 2-chlorophenyl, 3-nitrophenyl, 3-aminophenyl, 3-acetamidophenyl, 2-amino-3-(aminomethyl)phenyl, 6-

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methyl-3-acetamidophenyl, 6-methyl-2-aminophenyl, 6-methyl-2,3-diaminophenyl, 2-amino-3-methylphenyl, 4,6-dimethyl-2-aminophenyl, 4-hydroxyphenyl, 3-methyl-4-hydroxyphenyl, 4-(2-methoxyphenyl)phenyl, 2-amino-1-naphthyl, 2-naphthyl, 3-amino-2-naphthyl, 1-methyl-3-amino-2-naphthyl, 2,3-diamino-1-naphthyl, 4,8-dimethoxy-2-naphthyl and the like.

"Aralkyl" and "arylalkyl", alone or in combination, means an alkyl radical as defined above in which at least one hydrogen atom, preferably 1-2, is replaced by an aryl radical as defined above, such as benzyl, 1-, 2-phenylethyl, dibenzylmethyl, hydroxyphenylmethyl, methylphenylmethyl, diphenylmethyl, dichlorophenylmethyl, 4-methoxyphenylmethyl and the like.

"Aralkoxy", alone or in combination, means an alkoxy radical as defined above in which at least one hydrogen atom, preferably 1-2, is replaced by an aryl radical as defined above, such as benzyloxy, 1-, 2-phenylethoxy, dibenzylmethoxy, hydroxyphenylmethoxy, methylphenylmethoxy, dichlorophenylmethoxy, 4-methoxyphenylmethoxy and the like.

25

"Aralkoxycarbonyl", alone or in combination, means a radical of the type "R-O-C(O)-" wherein "R-O-" is an aralkoxy radical as defined above and "-C(O)-" is a carbonyl radical.

30

"Alkanoyl", alone or in combination, means a radical of the type "R-C(O)-" wherein "R" is an alkyl radical as defined above and "-C(O)-" is a carbonyl radical. Examples of such alkanoyl radicals include acetyl, trifluoroacetyl, hydroxyacetyl, propionyl, butyryl, valeryl, 4-methylvaleryl, and the like.

"Alkanoylamino", alone or in combination, means a radical of the type "R-C(O)-NH-" wherein "R-C(O)-" is an alkanoyl radical as defined above, wherein the amino radical may optionally be substituted, such as with
5 alkyl, aryl, aralkyl, cycloalkyl, cycloalkylalkyl and the like.

"Aminocarbonyl", alone or in combination, means an amino substituted carbonyl (carbamoyl) radical, wherein the
10 amino radical may optionally be mono- or di-substituted, such as with alkyl, aryl, aralkyl, cycloalkyl, cycloalkylalkyl, alkanoyl, alkoxycarbonyl, aralkoxycarbonyl and the like.

15 "Aminosulfonyl", alone or in combination, means an amino substituted sulfonyl radical.

"Benzo", alone or in combination, means the divalent radical $C_6H_4=$ derived from benzene. "Benzo fused" forms
20 a ring system in which benzene and a cycloalkyl or aryl group have two carbons in common, for example tetrahydronaphthylene and the like.

"Bicyclic" as used herein is intended to include both
25 fused ring systems, such as naphthyl and β -carbolinyl, and substituted ring systems, such as biphenyl, phenylpyridyl and diphenylpiperazinyl.

"Cycloalkyl", alone or in combination, means a saturated
30 or partially saturated, preferably one double bond, monocyclic, bicyclic or tricyclic carbocyclic alkyl radical, preferably monocyclic, containing preferably 5-12 carbon atoms (C_5-C_{12}), more preferably 5-10 carbon atoms (C_5-C_{10}), even more preferably 5-7 carbon atoms
35 (C_5-C_7), which is optionally benzo fused or heterocyclofused and which is optionally substituted as defined herein with respect to the definition of aryl. Examples

of such cycloalkyl radicals include cyclopentyl,
cyclohexyl, dihydroxycyclohexyl,
ethylenedioxcyclohexyl, cycloheptyl, octahydronaphthyl,
tetrahydronaphthyl, octahydroquinoliny1,
5 dimethoxytetrahydronaphthyl, 2,3-dihydro-1H-indenyl,
azabicyclo[3.2.1]octyl and the like.

"Heteroatoms" means nitrogen, oxygen and sulfur
heteroatoms.

10

"Heterocyclo fused" forms a ring system in which a
heterocyclyl or heteroaryl group of 5-6 ring members and
a cycloalkyl or aryl group have two carbons in common,
for example indole, isoquinoline, tetrahydroquinoline,
15 methylenedioxybenzene and the like.

"Heterocyclyl" means a saturated or partially
unsaturated, preferably one double bond, monocyclic or
bicyclic, preferably monocyclic, heterocycle radical
20 containing at least one, preferably 1 to 4, more
preferably 1 to 3, even more preferably 1-2, nitrogen,
oxygen or sulfur atom ring member and having preferably
3-8 ring members in each ring, more preferably 5-8 ring
members in each ring and even more preferably 5-6 ring
25 members in each ring. "Heterocyclyl" is intended to
include sulfone and sulfoxide derivatives of sulfur ring
members and N-oxides of tertiary nitrogen ring members,
and carbocyclic fused, preferably 3-6 ring carbon atoms
and more preferably 5-6 ring carbon atoms, and benzo
30 fused ring systems. "Heterocyclyl" radicals may
optionally be substituted on at least one, preferably 1-
4, more preferably 1-3, even more preferably 1-2, carbon
atoms by halogen, alkyl, alkoxy, hydroxy, oxo, thioxo,
aryl, aralkyl, heteroaryl, heteroaralkyl, amidino, N-
35 alkylamidino, alkoxycarbonylamino, alkylsulfonylamino
and the like, and/or on a secondary nitrogen atom by
hydroxy, alkyl, aralkoxycarbonyl, alkanoyl,

alkoxycarbonyl, heteroaralkyl, aryl or aralkyl radicals. More preferably, "heterocyclyl", alone or in combination, is a radical of a monocyclic or bicyclic saturated heterocyclic ring system having 5-8 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally partially unsaturated or benzo-fused and optionally substituted by 1-2 oxo or thioxo radicals. Examples of such heterocyclyl radicals include pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, thiamorpholinyl, 4-benzyl-piperazin-1-yl, pyrimidinyl, tetrahydrofuryl, pyrazolidonyl, pyrazolinyl, pyridazinonyl, pyrrolidonyl, tetrahydrothienyl and its sulfoxide and sulfone derivatives, 2,3-dihydroindolyl, tetrahydroquinolinyl, 1,2,3,4-tetrahydroisoquinolinyl, 1,2,3,4-tetrahydro-1-oxo-isoquinolinyl, 2,3-dihydrobenzofuryl, benzopyranyl, methylenedioxyphenyl, ethylenedioxyphenyl and the like.

"Heteroaryl" means a monocyclic or bicyclic, preferably monocyclic, aromatic heterocycle radical, having at least one, preferably 1 to 4, more preferably 1 to 3, even more preferably 1-2, nitrogen, oxygen or sulfur atom ring members and having preferably 5-6 ring members in each ring, which is optionally saturated carbocyclic fused, preferably 3-4 carbon atoms (C₃-C₄) to form 5-6 ring membered rings and which is optionally substituted as defined above with respect to the definitions of aryl. Examples of such heteroaryl groups include imidazolyl, 1-benzyloxycarbonylimidazol-4-yl, pyrrolyl, pyrazolyl, pyridyl, 3-(2-methyl)pyridyl, 3-(4-trifluoromethyl)pyridyl, pyrimidinyl, 5-(4-trifluoromethyl)pyrimidinyl, pyrazinyl, triazolyl, furyl, thienyl, oxazolyl, thiazolyl, indolyl, quinolinyl, 5,6,7,8-tetrahydroquinolyl, 5,6,7,8-tetrahydroisoquinolinyl, quinoxalinyl, benzothiazolyl, benzofuryl, benzimidazolyl, benzoxazolyl and the like.

"Heteroaralkyl" and "heteroarylalkyl," alone or in combination, means an alkyl radical as defined above in which at least one hydrogen atom, preferably 1-2, is
5 replaced by a heteroaryl radical as defined above, such as 3-furylpropyl, 2-pyrrolyl propyl, chloroquinolinylmethyl, 2-thienylethyl, pyridylmethyl, 1-imidazolethyl and the like.

10 "Halogen" and "halo", alone or in combination, means fluoro, chloro, bromo or iodo radicals.

"Haloalkyl", alone or in combination, means an alkyl radical as defined above in which at least one hydrogen
15 atom, preferably 1-3, is replaced by a halogen radical, more preferably fluoro or chloro radicals. Examples of such haloalkyl radicals include 1,1,1-trifluoroethyl, chloromethyl, 1-bromoethyl, fluoromethyl, difluoromethyl, trifluoromethyl,
20 bis(trifluoromethyl)methyl and the like.

"Pharmacologically acceptable salt" means a salt prepared by conventional means, and are well known by those skilled in the art. The "pharmacologically
25 acceptable salts" include basic salts of inorganic and organic acids, including but not limited to hydrochloric acid, hydrobromic acid, sulphuric acid, phosphoric acid, methanesulphonic acid, ethanesulfonic acid, malic acid, acetic acid, oxalic acid, tartaric acid, citric acid,
30 lactic acid, fumaric acid, succinic acid, maleic acid, salicylic acid, benzoic acid, phenylacetic acid, mandelic acid and the like. When compounds of the invention include an acidic function such as a carboxy group, then suitable pharmaceutically acceptable cation
35 pairs for the carboxy group are well known to those skilled in the art and include alkaline, alkaline earth, ammonium, quaternary ammonium cations and the like. For

additional examples of "pharmacologically acceptable salts," see *infra* and Berge et al, *J. Pharm. Sci.* **66**, 1 (1977).

5 "Cytokine" means a secreted protein that affects the functions of other cells, particularly as it relates to the modulation of interactions between cells of the immune system or cells involved in the inflammatory response. Examples of cytokines include but are not
10 limited to interleukin 1 (IL-1), preferably IL-1 β , interleukin 6 (IL-6), interleukin 8 (IL-8) and TNF, preferably TNF- α (tumor necrosis factor- α).

"TNF, IL-1, IL-6, and/or IL-8 mediated disease or
15 disease state" means all disease states wherein TNF, IL-1, IL-6, and/or IL-8 plays a role, either directly as TNF, IL-1, IL-6, and/or IL-8 itself, or by TNF, IL-1, IL-6, and/or IL-8 inducing another cytokine to be released. For example, a disease state in which IL-1
20 plays a major role, but in which the production of or action of IL-1 is a result of TNF, would be considered mediated by TNF.

"Leaving group" generally refers to groups readily
25 displaceable by a nucleophile, such as an amine, a thiol or an alcohol nucleophile. Such leaving groups are well known in the art. Examples of such leaving groups include, but are not limited to, N-hydroxysuccinimide, N-hydroxybenzotriazole, halides, triflates, tosylates
30 and the like. Preferred leaving groups are indicated herein where appropriate.

"Protecting group" generally refers to groups well known
in the art which are used to prevent selected reactive
35 groups, such as carboxy, amino, hydroxy, mercapto and the like, from undergoing undesired reactions, such as nucleophilic, electrophilic, oxidation, reduction and the

like. Preferred protecting groups are indicated herein where appropriate. Examples of amino protecting groups include, but are not limited to, aralkyl, substituted aralkyl, cycloalkenylalkyl and substituted cycloalkenyl alkyl, allyl, substituted allyl, acyl, alkoxycarbonyl, aralkoxycarbonyl, silyl and the like. Examples of aralkyl include, but are not limited to, benzyl, ortho-methylbenzyl, trityl and benzhydryl, which can be optionally substituted with halogen, alkyl, alkoxy, hydroxy, nitro, acylamino, acyl and the like, and salts, such as phosphonium and ammonium salts. Examples of aryl groups include phenyl, naphthyl, indanyl, anthracenyl, 9-(9-phenylfluorenyl), phenanthrenyl, durenyl and the like. Examples of cycloalkenylalkyl or substituted cycloalkylenylalkyl radicals, preferably have 6-10 carbon atoms, include, but are not limited to, cyclohexenyl methyl and the like. Suitable acyl, alkoxycarbonyl and aralkoxycarbonyl groups include benzyloxycarbonyl, t-butoxycarbonyl, iso-butoxycarbonyl, benzoyl, substituted benzoyl, butyryl, acetyl, tri-fluoroacetyl, tri-chloro acetyl, phthaloyl and the like. A mixture of protecting groups can be used to protect the same amino group, such as a primary amino group can be protected by both an aralkyl group and an aralkoxycarbonyl group. Amino protecting groups can also form a heterocyclic ring with the nitrogen to which they are attached, for example, 1,2-bis(methylene)benzene, phthalimidyl, succinimidyl, maleimidyl and the like and where these heterocyclic groups can further include adjoining aryl and cycloalkyl rings. In addition, the heterocyclic groups can be mono-, di- or tri-substituted, such as nitrophthalimidyl. Amino groups may also be protected against undesired reactions, such as oxidation, through the formation of an addition salt, such as hydrochloride, toluenesulfonic acid, trifluoroacetic acid and the like. Many of the amino protecting groups are also suitable for protecting carboxy, hydroxy and mercapto groups. For example,

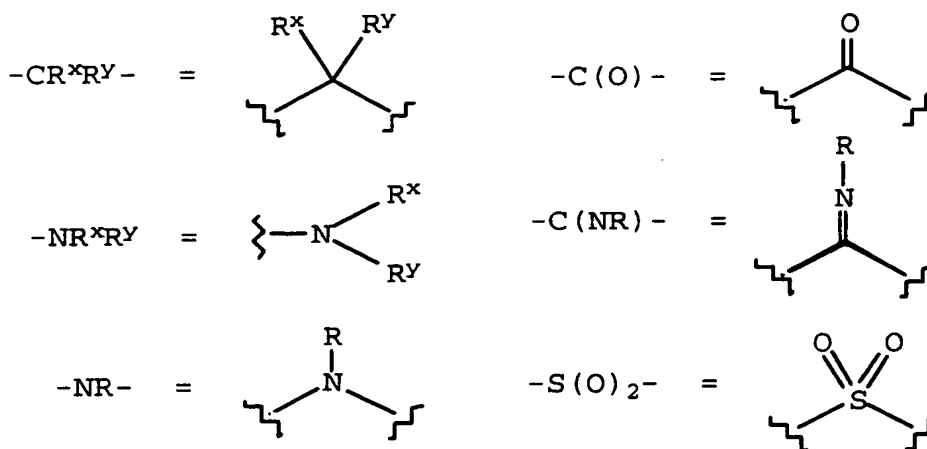
aralkyl groups. Alkyl groups are also suitable groups for protecting hydroxy and mercapto groups, such as tert-butyl.

Silyl protecting groups are silicon atoms
5 optionally substituted by one or more alkyl, aryl and aralkyl groups. Suitable silyl protecting groups include, but are not limited to, trimethylsilyl, triethylsilyl, tri-isopropylsilyl, tert-butyl-
10 dimethylsilyl, dimethylphenylsilyl, 1,2-bis(dimethylsilyl)benzene, 1,2-bis(dimethylsilyl)ethane and diphenylmethylsilyl. Silylation of an amino groups provide mono- or di-silylamino groups. Silylation of aminoalcohol compounds can lead to a N,N,O-tri-silyl derivative. Removal of the silyl function from a silyl
15 ether function is readily accomplished by treatment with, for example, a metal hydroxide or ammonium fluoride reagent, either as a discrete reaction step or in situ during a reaction with the alcohol group. Suitable silylating agents are, for example,
20 trimethylsilyl chloride, tert-butyl-dimethylsilyl chloride, phenyldimethylsilyl chloride, diphenylmethylsilyl chloride or their combination products with imidazole or DMF. Methods for silylation of amines and removal of silyl protecting groups are well known to
25 those skilled in the art. Methods of preparation of these amine derivatives from corresponding amino acids, amino acid amides or amino acid esters are also well known to those skilled in the art of organic chemistry including amino acid/amino acid ester or aminoalcohol
30 chemistry.

Protecting groups are removed under conditions which will not affect the remaining portion of the molecule. These methods are well known in the art and include acid hydrolysis, hydrogenolysis and the like. A
35 preferred method involves removal of a protecting group, such as removal of a benzyloxycarbonyl group by hydrogenolysis utilizing palladium on carbon in a

suitable solvent system such as an alcohol, acetic acid, and the like or mixtures thereof. A t-butoxycarbonyl protecting group can be removed utilizing an inorganic or organic acid, such as HCl or trifluoroacetic acid, in a suitable solvent system, such as dioxane or methylene chloride. The resulting amino salt can readily be neutralized to yield the free amine. Carboxy protecting group, such as methyl, ethyl, benzyl, tert-butyl, 4-methoxyphenylmethyl and the like, can be removed under hydrolysis and hydrogenolysis conditions well known to those skilled in the art.

The symbols used above have the following meanings:



Procedures for preparing the compounds of this invention are set forth below. It should be noted that the general procedures are shown as it relates to preparation of compounds having unspecified stereochemistry. However, such procedures are generally applicable to those compounds of a specific stereochemistry, e.g., where the stereochemistry about a group is (S) or (R). In addition, the compounds having one stereochemistry (e.g., (R)) can often be utilized to produce those having opposite stereochemistry (i.e., (S)) using well-known methods, for example, by inversion.

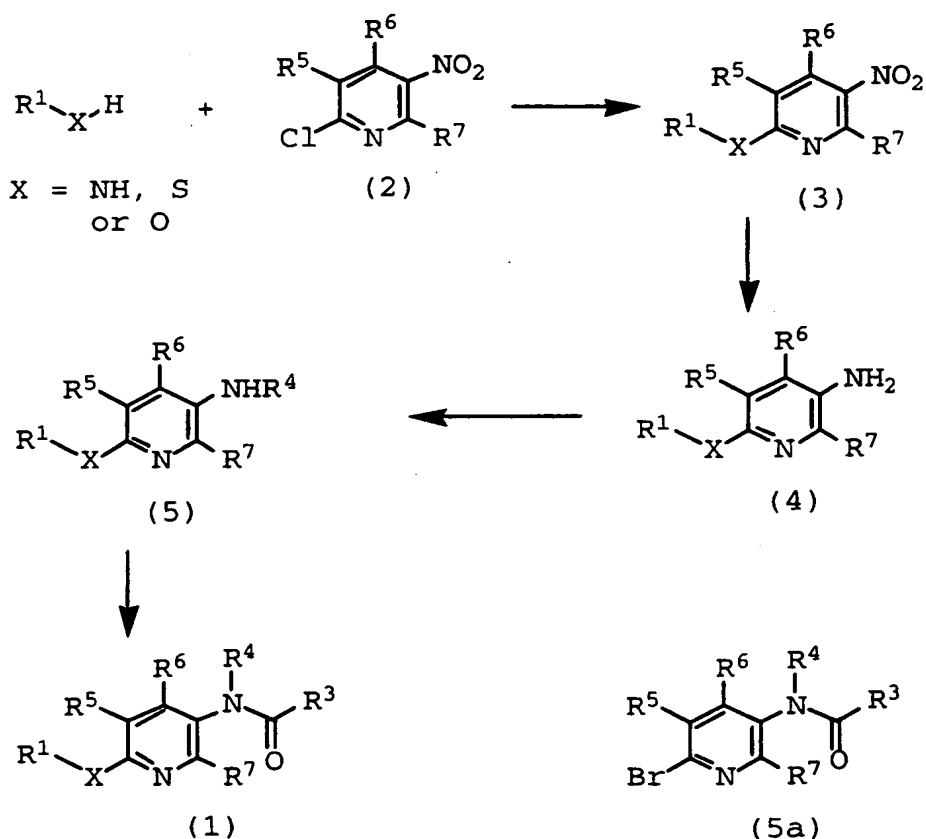
Preparation of Compounds of Formula I

The compounds of the present invention represented
5 by Formula I above can be prepared utilizing the
following general procedures. Hetero-aromatic Nitrogen
Compounds; Pyrroles and Pyridines: Schofield, Kenneth;
Plenum Press, New York, NY; (1967) and Advances in
Nitrogen Heterocycles: JAI Press, Greenwich, CN; (1995)
10 describe procedures and references that may be useful in
preparing compounds of the present invention.

2-Halo-5-nitro-pyridine analogs (2) can be treated
with the appropriate amine, alcohol, phenol, or thiol
(R^1-X-H) in the presence of base or Cu(I) in an
15 appropriate solvent, such as THF, DMF, DME, DMSO and the
like, at a temperature from -20°C to 120°C to form 2-
substituted-5-nitropyridines (3) (Scheme I). Reduction
of the nitro group can be performed by treatment of (3)
with hydrogen gas in the presence of palladium on carbon
20 or Raney nickel, or alternatively, by treatment with
 SnCl_2 in an alcoholic solvent and in the presence or
absence of HCl to obtain 2-substituted-5-aminopyridines
(4). The aminopyridines (4) may be alkylated using
alkylhalides and an appropriate base or by reductive
25 alkylation employing the appropriate aldehyde or ketone
in the presence of a reducing agent, such as sodium
triacetoxo borohydride, borane-THF and the like, to form
the substituted aminopyridines (5). Either (4) or (5)
may be acylated with an appropriate acid halide (e.g.,
30 $R^3\text{C}(\text{O})\text{Cl}$ or $R^3\text{C}(\text{O})\text{Br}$) in the presence of a base, such as
pyridine, DMAP and the like, or alternatively may be
acylated with an anhydride, either mixed or symmetrical,
or alternatively may be acylated by treatment with the
appropriate acid ($R^3\text{CO}_2\text{H}$) in the presence of a coupling
35 agent such as a carbodiimide reagent to form the final
product (1). Alternatively, substituted 2-bromo-5-
nitropyridine analogs may be reduced to, substituted 2-

bromo-5-aminopyridine analogs by the action of SnBr_2 in methanolic solvent. Subsequent acylation with an appropriate activated ester (i.e.: $\text{R}^3\text{CO}_2\text{H}$ in the presence of diisopropylcarbodiimide in methylene chloride as solvent) produces 2-bromopyridine-5-carboxamide compounds of structure (5a). Coupling of (5a) with an appropriate phenol in the presence of $\text{Cu}(\text{Ac})_2$ and K_2CO_3 in DMF at 140°C provides compounds of formula (1) where $\text{X} = \text{O}$.

10

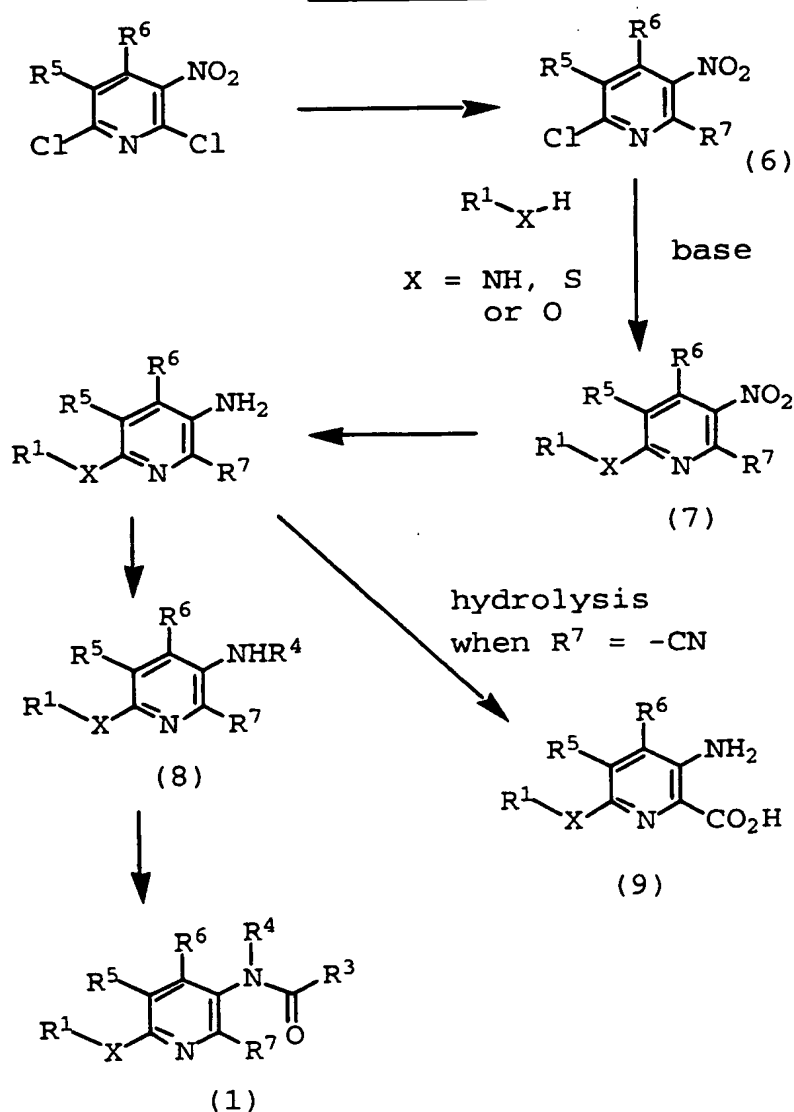
SCHEME I

6-Substituted-2-halo-5-nitropyridine analogs (6) may be prepared from 2,6-dichloro-5-nitropyridine according to the methods outlined in Scheme II. Treatment with one equivalent of an appropriate nucleophile of R^7 or a precursor thereof (such as, HO^- , RO^- , AcS^- , NC^- , RS^- and the like) provides (6). Subsequent reaction to form (7) (treatment with $\text{R}^1\text{-X-H}$

in the presence of base or Cu(I) in an appropriate solvent, such as THF, DMF, DME, DMSO and the like, at a temperature from -20°C to 120°C) and (8) (reduction of the nitro group and substitution with R⁴) is as

- 5 described in Scheme I (cf. Colbry, N.L. et al.; J. Heterocyclic Chem., 21: 1521-1525 (1984); Matsumoto, Jun-ichi, et al.; J. Heterocyclic Chem., 21: 673-679 (1984)). (8) may be reacted with an acid halide or an activated

10

SCHEME II

ester as shown in Scheme I to provide compounds of formula (1). Where R⁷ = CN, compounds of formula (8)

may be hydrolyzed to acids ($R' = CO_2H$) of formula (9) using acidic media such as HBr and the like. Utilizing the appropriate N-protecting groups, acids of formula (9) may be transformed into esters, amides and alcohols.

5 Compounds of formula (9) and derivatives described above may be reacted with an acid halide or an activated ester as shown in Scheme I to provide compounds of formula (1). Compounds of formula (8), where $R' = -CN$, may be reduced to the primary amine ($R' = -CH_2NH_2$) using

10 reagents such as BH_3 or hydrogen gas in the presence of palladium on carbon or Raney nickel. Subsequent manipulation and reaction of the primary amine may be performed in the presence of the pyridine-5-amine substituent due to its greater reactivity.

15 Specifically, compounds of formula (8) where $R' = -CH_2NH_2$ may be alkylated by treatment with an appropriate aldehyde or ketone in the presence of a reducing agent, such as sodium triacetoxy borohydride, or may be acylated by treatment with an appropriate activated

20 ester, chloroformate, isocyanate and the like, or may be sulfonylated by treatment with an appropriate sulfonyl halide. Alternatively, substituted 3-aminopyridine intermediates may be prepared from the corresponding nicotinamide compound using Hofmann's reaction.

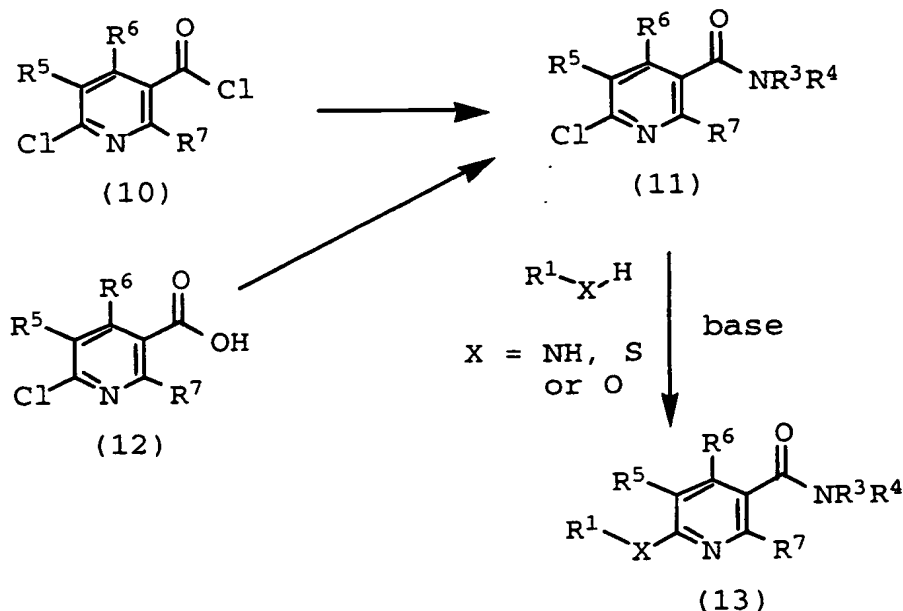
25 When R^6 and/or R^7 is an alkyl group, such as methyl, in compound (7), containing the appropriate protecting groups of or avoiding the presence of base sensitive groups, can be treated with strong base such as $NaNH_2$, $PhLi$, NaH or the like at temperatures from $-78^\circ C$ to $22^\circ C$

30 then treated with electrophiles, such as alkyl halides, aldehydes, ketones and the like (cf. Fuerst, Feustel; CHEMTECH; 10: 693-699 (1958); Nishigaki, S. et al.; Chem. Pharm. Bull.; 17: 1827-1831 (1969); Kaiser, Edwin M.; Tetrahedron; 39: 2055-2064 (1983)). Alternatively,

35 the alkyl group may be halogenated and the haloalkyl group may be reacted with a nucleophile, such as an amino group, alkoxy, alkylthiol and the like.

6-Chloronicotinoyl chloride analogs (10) are treated with the appropriate amine (R^3R^4NH) in the presence of base in an appropriate solvent, such as dichloromethane, acetonitrile, DMF, THF and the like, at a temperature from -20°C to 120°C to form nicotinamides (11) as shown in Scheme III. Alternatively, 6-chloronicotinic acid analogs (12) may be coupled with the appropriate amine via an anhydride, either mixed or symmetrical, or alternatively by treatment with the appropriate amine in the presence of a coupling agent such as a carbodiimide reagent to form the amide (11). 6-Chloronicotinamide analogs (11) are treated with the appropriate R^1-X-H in the presence of absence of base, or Cu(I) in an appropriate solvent, such as pyridine, ethylene glycol, DMF, DME, DMSO and the like, at a temperature from -20°C to 180°C to form the final product (13).

SCHEME III



20

Substituted halopyridines may be readily prepared from the corresponding pyridones using phosphorus oxychloride or pentachloride.

Amines of formula NHR^1R^2 and NHR^3R^4 are commercially available or can be readily prepared by those skilled in the art from commercially available starting materials. For example, an amide, nitro or cyano group can be

5 reduced under reducing conditions, such as in the presence of a reducing agent like lithium aluminum hydride and the like, to form the corresponding amine. Alkylation and acylation of amino groups are well known in the art. Chiral and achiral substituted amines can

10 be prepared from chiral amino acids and amino acid amides (for example, alkyl, aryl, heteroaryl, cycloalkyl, arylalkyl, heteroarylalkyl, cycloalkylalkyl and the like) using methods well known in the art, such as H. Brunner, P. Hankofer, U. Holzinger, B. Treitinger

15 and H. Schoenenberger, Eur. J. Med. Chem. 25, 35-44, 1990; M. Freiburger and R. B. Hasbrouck, J. Am. Chem. Soc. 82, 696-698, 1960; Dornow and Fust, Chem. Ber. 87, 984, 1954; M. Kojima and J. Fujita, Bull. Chem. Soc. Jpn. 55, 1454-1459, 1982; W. Wheeler and D. O'Bannon,

20 Journal of Labelled Compounds and Radiopharmaceuticals XXXI, 306, 1992; and S. Davies, N. Garrido, O. Ichihara and I. Walters, J. Chem. Soc., Chem. Commun. 1153, 1993.

Alkyl sulfonic acids, aryl sulfonic acids, heterocyclyl sulfonic acids, heteroaryl sulfonic acids,

25 alkylmercaptans, arylmercaptans, heterocyclylmercaptans, heteroarylmercaptans, alkylhalides, arylhalides, heterocyclylhalides, heteroarylhalides, and the like are commercially available or can be readily prepared from starting materials commercially available using standard

30 methods well known in the art.

Thioether derivatives can be converted into the corresponding sulfone or sulfoxide by oxidizing the thioether derivative with a suitable oxidation agent in a suitable solvent. Suitable oxidation agents include,

35 for example, hydrogen peroxide, sodium meta-perborate, oxone (potassium peroxy monosulfate), meta-chloroperoxybenzoic acid, periodic acid and the like,

including mixtures thereof. Suitable solvents include acetic acid (for sodium meta-perborate) and, for other peracids, ethers such as THF and dioxane, and acetonitrile, DMF and the like, including mixtures thereof.

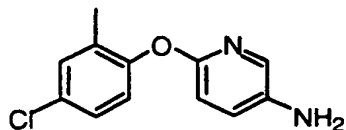
The chemical reactions described above are generally disclosed in terms of their broadest application to the preparation of the compounds of this invention. Occasionally, the reactions may not be applicable as described to each compound included within the disclosed scope. The compounds for which this occurs will be readily recognized by those skilled in the art. In all such cases, either the reactions can be successfully performed by conventional modifications known to those skilled in the art, e.g., by appropriate protection of interfering groups, by changing to alternative conventional reagents, by routine modification of reaction conditions, and the like, or other reactions disclosed herein or otherwise conventional, will be applicable to the preparation of the corresponding compounds of this invention. In all preparative methods, all starting materials are known or readily prepared from known starting materials.

Prodrugs of the compounds of this invention are also contemplated by this invention. A prodrug is an active or inactive compound that is modified chemically through in vivo physiological action, such as hydrolysis, metabolism and the like, into a compound of this invention following administration of the prodrug to a patient. The suitability and techniques involved in making and using prodrugs are well known by those skilled in the art. For a general discussion of prodrugs involving esters see Svensson and Tunek Drug Metabolism Reviews 165 (1988) and Bundgaard Design of Prodrugs, Elsevier (1985). Examples of a masked carboxylate anion include a variety of esters, such as alkyl (for example, methyl, ethyl), cycloalkyl (for

example, cyclohexyl), aralkyl (for example, benzyl, p-methoxybenzyl), and alkylcarbonyloxyalkyl (for example, pivaloyloxymethyl). Amines have been masked as
5 arylcarbonyloxymethyl substituted derivatives which are cleaved by esterases in vivo releasing the free drug and formaldehyde (Bunggaard J. Med. Chem. 2503 (1989)). Also, drugs containing an acidic NH group, such as imidazole, imide, indole and the like, have been masked with N-acyloxymethyl groups (Bundgaard Design of
10 Prodrugs, Elsevier (1985)). Hydroxy groups have been masked as esters and ethers. EP 039,051 (Sloan and Little, 4/11/81) discloses Mannich-base hydroxamic acid prodrugs, their preparation and use.

Without further elaboration, it is believed that
15 one skilled in the art can, using the preceding description, utilize the present invention to its fullest extent. The following preferred specific embodiments are, therefore, to be construed as merely illustrative, and not limitative of the remainder of the disclosure in any way whatsoever. The following
20 Examples illustrate the preparation of compounds of the present invention and intermediates useful in preparing the compounds of the present invention.

25

Example 1Preparation of 2-(4-Chloro-2-methyl-phenoxy)-5-amino-pyridine

30

Step A: 2-(4-Chloro-2-methyl-phenoxy)-5-nitropyridine

4-Chloro-2-methylphenol (101 mg, 0.71 mmol) was dissolved in tetrahydrofuran (2.1 mL) and the solution was treated with sodium hydride (60% dispersed in
35 mineral oil, 31 mg, 0.78 mmol). After stirring for 30

minutes at 22°C, 2-chloro-5-nitropyridine (101 mg, 0.64 mmol) was added and the reaction mixture was heated to reflux for 1 hour. The solution was cooled to ambient temperature, quenched with saturated aqueous NH_4Cl and concentrated in vacuo. The residue was redissolved in ethyl acetate then washed 2x with saturated NaHCO_3 , saturated NaCl , dried over anhydrous Na_2SO_4 and concentrated in vacuo.

10 Step B: 2-(4-Chloro-2-methyl-phenoxy)-5-amino-pyridine
2-(4-chloro-2-methyl-phenoxy)-5-nitropyridine (203 mg, 0.77 mmol) was dissolved in 95% ethanol (3 mL) and treated with 20% palladium hydroxide on carbon (50 mg). The reaction mixture was shaken in a hydrogen atmosphere (40 psi) for 1 hour. The solution was filtered through celite and concentrated in vacuo. MS (m/z): 234/236 (M+H)⁺; $\text{C}_{12}\text{H}_{11}\text{N}_2\text{OCl}$ requires 234.7.

Example 2

20 The compounds listed in Table 1 were prepared from 2-chloro-5-nitropyridine and the appropriate alcohol, amine or thiol in the same manner as 2-(4-Chloro-2-methyl-phenoxy)-5-amino-pyridine was prepared.

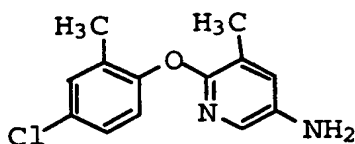
25

Table 1

| | MS (m/z) |
|---|-------------|
| 2-(4-Chloro-2-methylphenoxy)-5-amino-pyridine | 235 |
| 2-(4-Chloro-2,6-dimethylphenoxy)-5-amino-pyridine | 249 |
| 2-(2-Methyl-pyridin-3-yloxy)-5-amino-pyridine | 201 |
| 2-(4-Fluoro-2-methylphenoxy)-5-amino-pyridine | 218 |
| 2-(2-Isopropylphenoxy)-5-amino-pyridine | 228 |
| 2-(1-Naphthyloxy)-5-amino-pyridine | 236 |

| | |
|---|-----|
| 2-(Cyclohexyloxy)-5-amino-pyridine | 192 |
| 2-(2-Methylphenoxy)-5-amino-pyridine | 200 |
| 2-(2,4-Dimethylphenoxy)-5-amino-pyridine | 214 |
| 2-(4-Chlorophenoxy)-5-amino-pyridine | 222 |
| 2-(Phenoxy)-5-amino-pyridine | 186 |
| 2-(2-Methylcyclohexylamino)-5-amino-pyridine | 205 |
| 2-(Cyclohexylamino)-5-amino-pyridine | 191 |
| 2-(2-Methylanilino)-5-amino-pyridine | 199 |
| 2-(4-Chloro-2-methylanilino)-5-amino-pyridine | 233 |
| 2-(2,4-Dimethylanilino)-5-amino-pyridine | 212 |
| 2-(4-Chloro-2-methylthiophenoxy)-5-amino-pyridine | 251 |

Example 3



5 Preparation of 2-(4-Chloro-2-methyl-phenoxy)-3-methyl-5-
amino-pyridine

Step A: 2-(4-Chloro-2-methyl-phenoxy)-3-methyl-5-
nitropyridine

10 Sodium hydride (60% in mineral oil, 1.08 g, 27 mmol) was washed 3x with hexanes then a solution of 4-chloro-2-methylphenol (3.50 g, 24.5 mmol) dissolved in tetrahydrofuran (40 mL) was added. The solution was stirred for 20 minutes then 2-chloro-3-methyl-5-nitropyridine (4.02 g, 23.3 mmol) was added and the
15 reaction mixture was heated to reflux for 3 hours. After cooling, the mixture was concentrated in *vacuo* then dissolved in ethyl acetate and washed with water, 3x with saturated NaHCO₃ and saturated NaCl then dried over Na₂SO₄ and concentrated in *vacuo*.

Step B: 2-(4-Chloro-2-methyl-phenoxy)-3-methyl-5-amino-pyridine

2-(4-chloro-2-methyl-phenoxy)-3-methyl-5-nitropyridine
5 (5.8 g, 20.8 mmol) was dissolved in 95% ethanol (50 mL) and treated with 20% palladium hydroxide on carbon (350 mg). The reaction mixture was shaken in a hydrogen atmosphere (40 psi) for 1 hour. The solution was filtered through celite and concentrated in vacuo
10 followed by chromatography on SiO₂ using 1:1 ethyl acetate / hexanes as eluant. MS (m/z): 248/250 (M+H)⁺; C₁₃H₁₃N₂OCl requires 248.7.

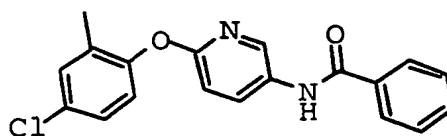
Example 4

15 The compounds listed in Table 2 were prepared from substituted 2-chloro-5-nitropyridine and 4-chloro-2-methylphenol in the same manner as 2-(4-Chloro-2-methyl-phenoxy)-3-methyl-5-amino-pyridine was prepared.

20 Table 2

| | MS (m/z) |
|---|-------------|
| 2-(4-Chloro-2-methyl-phenoxy)-4-methyl-5-amino-pyridine | 249 |
| 6-(4-Chloro-2-methyl-phenoxy)-2-methyl-3-amino-pyridine | 249 |
| 6-(4-Chloro-2-methyl-phenoxy)-2,3-diamino-pyridine | 250 |

Example 5



Preparation of N-(2-(4-Chloro-2-methyl-phenoxy)-pyridin-
25 5-yl)-benzamide

2-(4-Chloro-2-methyl-phenoxy)-5-aminopyridine (211 mg, 0.90 mmol) was dissolved in methylene chloride (2.7 mL) then treated with triethylamine (0.19 mL, 1.35 mmol) followed by benzoyl chloride (0.13 mL, 1.12 mmol). The reaction mixture was stirred for 3 hours at 22°C then saturated aqueous NaHCO₃ was added and the mixture was stirred for another hour. The organic layer was separated and washed 2x with 6% aqueous NaHCO₃, dried over Na₂SO₄ and concentrated in vacuo. The residue was chromatographed on silica gel using 1:1 ethyl acetate / hexane as eluent. The product was recovered as a white solid. MS (m/z): 338/340 (M+H)⁺; C₁₉H₁₅N₂O₂Cl requires 338.8.

Example 6

The compounds listed in Table 3 were prepared from substituted 5-aminopyridine compounds and the appropriate acid chloride in the same manner as N-(2-(4-Chloro-2-methyl-phenoxy)-pyridin-5-yl)-benzamide was prepared.

Table 3

| | MS (m/z) |
|---|-------------|
| 2-(4-Chloro-2-methyl-phenoxy)-5-(3-pyridylcarbonylamino)pyridine | 340 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 408 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-(4-pyridylcarbonylamino)pyridine | 340 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((4-methoxyphenyl)carbonylamino)pyridine | 369 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((4-pentylphenyl)carbonylamino)pyridine | 409 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-(2-naphthylcarbonylamino)pyridine | 389 |

| | |
|--|-----|
| 2-(4-Chloro-2-methyl-phenoxy)-5-(2-thienylcarbonylamino)pyridine | 345 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((3,5-dimethyl-4-isoxazolyl)carbonylamino)pyridine | 358 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((5-benzo[1,3]dioxol-yl)carbonylamino)pyridine | 383 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((5-tert-butyl-2-methyl-2H-pyrazol-3-yl)carbonylamino)pyridine | 399 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-benzo[b]thiophenyl)carbonylamino)pyridine | 395 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-methoxyphenyl)carbonylamino)pyridine | 369 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((3,5-dichlorophenyl)carbonylamino)pyridine | 408 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 367 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-methylphenyl)carbonylamino)pyridine | 353 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-nitrophenyl)carbonylamino)pyridine | 384 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-acetoxyphenyl)carbonylamino)pyridine | 397 |
| 2-(4-chloro-2,6-dimethylphenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 422 |
| 2-(4-chloro-2,6-dimethylphenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 381 |
| 2-(2-methyl-pyridin-3-yloxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 374 |
| 2-(2-methyl-pyridin-3-yloxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 333 |
| 2-(2-methyl-pyridin-3-yloxy)-5-((2-methylphenyl)carbonylamino)pyridine | 319 |

| | |
|---|-----|
| 2-(4-fluoro-2-methylphenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 391 |
| 2-(4-fluoro-2-methylphenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 350 |
| 2-(4-fluoro-2-methylphenoxy)-5-((2-methylphenyl)carbonylamino)pyridine | 336 |
| 2-(4-fluoro-2-methylphenoxy)-5-((2-trifluoromethylphenyl)carbonylamino)pyridine | 390 |
| 2-(4-fluoro-2-methylphenoxy)-5-((2-fluorophenyl)carbonylamino)pyridine | 340 |
| 2-(2-isopropylphenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 401 |
| 2-(2-isopropylphenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 360 |
| 2-(2-isopropylphenoxy)-5-((2-methylphenyl)carbonylamino)pyridine | 346 |
| 2-(1-naphthyloxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 409 |
| 2-(1-naphthyloxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 368 |
| 2-(1-naphthyloxy)-5-((2-methylphenyl)carbonylamino)pyridine | 354 |
| 2-(cyclohexyloxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 365 |
| 2-(cyclohexyloxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 324 |
| 2-(cyclohexyloxy)-5-((2-chlorophenyl)carbonylamino)pyridine | 331 |
| 2-(cyclohexyloxy)-5-((2-methylphenyl)carbonylamino)pyridine | 310 |
| 2-(2-methylphenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 373 |

| | |
|---|-----|
| 2-(2-methylphenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 332 |
| 2-(2-methylphenoxy)-5-((2-chlorophenyl)carbonylamino)pyridine | 339 |
| 2-(2-methylphenoxy)-5-((2-methylphenyl)carbonylamino)pyridine | 318 |
| 2-(2,4-dimethylphenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 387 |
| 2-(2,4-dimethylphenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 346 |
| 2-(2,4-dimethylphenoxy)-5-((2-chlorophenyl)carbonylamino)pyridine | 353 |
| 2-(2,4-dimethylphenoxy)-5-((2-methylphenyl)carbonylamino)pyridine | 332 |
| 2-(4-chlorophenoxy)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 394 |
| 2-(4-chlorophenoxy)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 353 |
| 2-(2-methylcyclohexylamino)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 378 |
| 2-(2-methylcyclohexylamino)-5-((2-methylphenyl)carbonylamino)pyridine | 323 |
| 2-(cyclohexylamino)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 364 |
| 2-(cyclohexylamino)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 323 |
| 2-(cyclohexylamino)-5-((2-methylphenyl)carbonylamino)pyridine | 309 |
| 2-(2-methylanilino)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 372 |
| 2-(2-methylanilino)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 331 |

| | |
|--|-----|
| 2-(2-methylanilino)-5-((2-methylphenyl)carbonylamino)pyridine | 317 |
| 2-(4-chloro-2-methylanilino)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 407 |
| 2-(4-chloro-2-methylanilino)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 366 |
| 2-(4-chloro-2-methylanilino)-5-((2-methylphenyl)carbonylamino)pyridine | 352 |
| 2-(2,4-dimethylanilino)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 386 |
| 2-(2,4-dimethylanilino)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 345 |
| 2-(2,4-dimethylanilino)-5-((2-methylphenyl)carbonylamino)pyridine | 331 |
| 2-(2,4-dimethylanilino)-5-((2-chlorophenyl)carbonylamino)pyridine | 352 |
| 2-(2,4-dimethylanilino)-5-((2-fluorophenyl)carbonylamino)pyridine | 335 |
| 2-(4-chloro-2-methyl-thiophenyl)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 424 |
| 2-(4-chloro-2-methyl-thiophenyl)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 383 |
| 2-(4-chloro-2-methyl-thiophenyl)-5-((2-methylphenyl)carbonylamino)pyridine | 369 |

Example 7

The compounds listed in Table 4 were prepared from substituted 5-aminopyridine compounds and the appropriate acid chloride in the same manner as N-(2-(4-Chloro-2-methyl-phenoxy)-pyridin-5-yl)-benzamide was prepared.

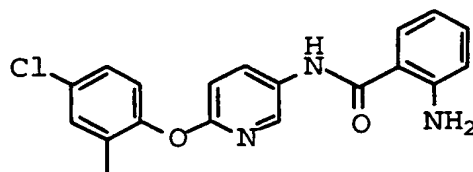
Table 4

| | MS (m/z) |
|---|-------------|
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dichlorophenyl)carbonylamino)-3-methyl-pyridine | 422 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-chlorophenyl)carbonylamino)-3-methyl-pyridine | 387 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-methylphenyl)carbonylamino)-3-methyl-pyridine | 367 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dimethylphenyl)carbonylamino)-3-methyl-pyridine | 332 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dichlorophenyl)carbonylamino)-4-methyl-pyridine | 422 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-fluoro-6-trifluoromethylphenyl)carbonylamino)-4-methyl-pyridine | 439 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,4,6-triisopropylphenyl)carbonylamino)-4-methyl-pyridine | 479 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-methylphenyl)carbonylamino)-6-methyl-pyridine | 367 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-chlorophenyl)carbonylamino)-6-methyl-pyridine | 387 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dichlorophenyl)carbonylamino)-6-methyl-pyridine | 422 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dichlorophenyl)carbonylamino)-6-amino-pyridine | 423 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2-chlorophenyl)carbonylamino)-6-amino-pyridine | 388 |
| 2-(4-Chloro-2-methyl-phenoxy)-5-((2,6-dimethylphenyl)carbonylamino)-6-amino-pyridine | 382 |

2-(4-Chloro-2-methyl-phenoxy)-5-((2-methylphenyl)carbonylamino)-6-amino-pyridine

368

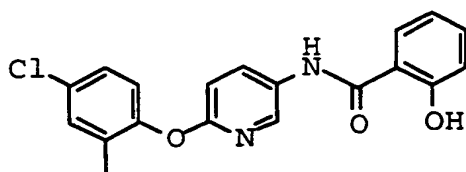
Example 8



5 Preparation of 2-Amino-N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-benzamide

N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-2-nitrobenzamide (301 mg, 0.7 mmol) was dissolved in 95% ethanol (4 mL) and treated with 20% palladium hydroxide
10 on carbon (Pearlman's catalyst, 50 mg) and subjected to a hydrogen atmosphere (40 psi) for 2 hours. The catalyst was removed by filtration and the solvents were removed in vacuo. The product was purified by chromatography on SiO₂ using 1:1 ethyl acetate / hexanes
15 as eluent. MS (m/z): 353/355 (M+H)⁺; C₁₉H₁₆N₃O₂Cl requires 353.8.

Example 9

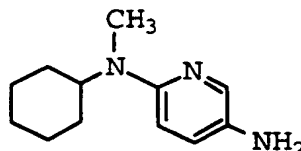


20 Preparation of N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-2-hydroxy-benzamide

Acetic acid 2-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-ylcarbonyl)-phenyl ester (304 mg, 0.77 mmol) dissolved in tetrahydrofuran (3.8 mL) was treated with an aqueous
25 lithium hydroxide solution (1.0 M, 3.8 mL, 3.8 mmol). The solution was stirred for 30 minutes at 22°C then quenched with aqueous saturated NH₄Cl. The mixture was diluted with ethyl acetate then the organics were washed

with water, 2x saturated NaHCO_3 , saturated NaCl , dried over Na_2SO_4 and concentrated in vacuo. MS (m/z): 354/356 ($\text{M}+\text{H}$) $^+$; $\text{C}_{19}\text{H}_{15}\text{N}_2\text{O}_3\text{Cl}$ requires 354.8.

5

Example 10Preparation of 2-(N-Cyclohexyl-N-methylamino)-5-amino-pyridine10 Step A: 2-(Cyclohexylamino)-5-nitro-pyridine

Sodium hydride (60% dispersion in mineral oil, 1.99 g, 49.8 mmol) was washed 3x with hexanes then a solution of cyclohexylamine (3.8 mL, 33.2 mmol) dissolved in tetrahydrofuran (50 mL) was added. After stirring for 30 minutes at 22 °C, 2-chloro-5-nitropyridine (5.00 g, 31.5 mmol) was added and the reaction mixture was heated to reflux for 3 hours. The solution was cooled to ambient temperature, quenched with saturated aqueous NH_4Cl and concentrated in vacuo. The residue was redissolved in ethyl acetate then washed 2x with saturated NaHCO_3 , saturated NaCl , dried over anhydrous Na_2SO_4 and concentrated in vacuo. The product was recovered as a brown oil.

25 Step B: 2-(N-Cyclohexyl-N-methylamino)-5-nitro-pyridine

Sodium hydride (60% dispersion in mineral oil, 0.38 g, 9.48 mmol) was washed 3x with hexanes then a solution of 2-cyclohexylamino-5-nitropyridine (1.88 g, 8.5 mmol) dissolved in dimethylformamide (20 mL) was added. After stirring for 30 minutes at 22°C, the reaction mixture was cooled to 0°C and methyl iodide (0.55 mL, 8.9 mmol) was added. The solution was stirred for 1.5 hours at 0°C followed by quenching with saturated aqueous NH_4Cl . The reaction mixture was diluted with ethyl acetate and

extracted 5x with water (200 mL), saturated NaCl, dried over Na₂SO₄ and concentrated in vacuo oil was chromatographed on SiO₂ using 2:1 hexanes / ethyl acetate as eluent.

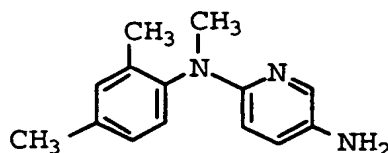
5

Step C: 2-(N-Cyclohexyl-N-methylamino)-5-amino-pyridine
Cyclohexyl-methyl-(5-nitro-pyridin-2-yl)-amine (1.72 g, 7.3 mmol) was dissolved in ethanol (80 mL) and treated with 20% palladium hydroxide on carbon (Pearlman's catalyst, 0.5 g) and the mixture was shaken under a hydrogen atmosphere (50 psi) for 6 hours. The catalyst was removed by filtration through celite then the filtrate was concentrated in vacuo and the resultant oil was chromatographed on SiO₂ using 1:1 ethyl acetate / hexanes as eluent. MS (m/z): 206 (M+H)⁺; C₁₂H₁₉N₃ requires 205.3.

10

15

Example 11

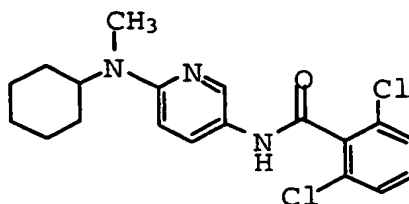


Preparation of 2-(N-(2,4-dimethylphenyl)-N-methylamino)-5-amino-pyridine
2-(N-(2,4-dimethylphenyl)-N-methylamino)-5-amino-pyridine was prepared from 1-amino-2,4-dimethylbenzene and 2-chloro-5-nitropyridine in the same manner as 2-(N-Cyclohexyl-N-methylamino)-5-amino-pyridine was prepared.

20

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Example 12



Preparation of 2,6-Dichloro-N-(2-(N'-cyclohexyl-N'-methylamino)-pyridin-5-yl)-benzamide

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2-(N-Cyclohexyl-N-methylamino)-5-amino-pyridine (26 mg, 0.13 mmol) dissolved in methylene chloride (0.25 mL) was treated with triethylamine (0.026 mL, 0.18 mmol) followed by a solution of 2,6-dichlorobenzoyl chloride (31 mg, 0.15 mmol) dissolved in methylene chloride (0.15 mL). The reaction mixture was shaken at 22°C for 18 hours followed by quenching with saturated aqueous NH_4Cl and stirring for an additional 5 hours. The organic layer was separated and dried over Na_2SO_4 , then concentrated *in vacuo*. The crude product was purified by chromatography on SiO_2 using 1:1 ethyl acetate / hexane as eluent. MS (m/z): 378/380 ($\text{M}+\text{H}$)⁺; $\text{C}_{19}\text{H}_{21}\text{N}_3\text{OCl}$ requires 377.

15

Example 13

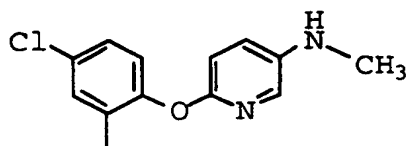
The compounds listed in Table 5 were prepared from substituted 5-aminopyridine compounds and the appropriate acid chloride in the same manner as 2,6-Dichloro-N-(2-(N'-cyclohexyl-N'-methylamino)-pyridin-5-yl)-benzamide was prepared.

Table 5

| | MS (m/z) |
|--|-----------------|
| 2-(N-cyclohexyl-N-methylamino)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 378 |
| 2-(N-cyclohexyl-N-methylamino)-5-((2-chlorophenyl)carbonylamino)pyridine | 344 |
| 2-(N-cyclohexyl-N-methylamino)-5-((2-methylphenyl)carbonylamino)pyridine | 323 |
| 2-(N-cyclohexyl-N-methylamino)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 337 |
| 2-(2,4-dimethylphenyl)-5-((2,6-dimethylphenyl)carbonylamino)pyridine | 359 |

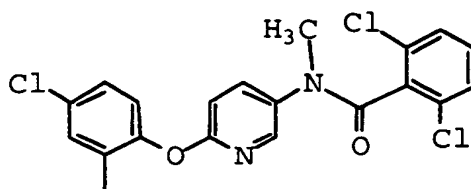
| | |
|--|-----|
| 2-(2,4-dimethylphenyl)-5-((2-methylphenyl)carbonylamino)pyridine | 345 |
| 2-(2,4-dimethylphenyl)-5-((2-chlorophenyl)carbonylamino)pyridine | 366 |
| 2-(2,4-dimethylphenyl)-5-((2-fluorophenyl)carbonylamino)pyridine | 349 |
| 2-(2,4-dimethylphenyl)-5-((2,6-dichlorophenyl)carbonylamino)pyridine | 400 |

Example 14

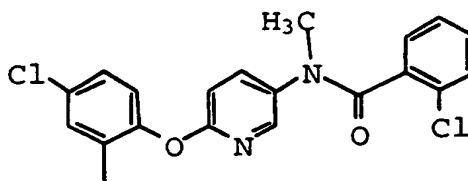


5 Preparation of 2-(4-Chloro-2-methyl-phenoxy)-5-(N-methylamino)pyridine

2-(4-Chloro-2-methyl-phenoxy)-5-aminopyridine (2.15 g, 9.16 mmol) was combined with powdered sodium hydroxide (1.46 g, 36.6 mmol), potassium carbonate (1.27 g, 9.16 mmol), tetrabutyl ammonium bromide (60 mg, 0.18 mmol) and toluene (10 mL) was stirred for 1 hour at 35°C. A solution of dimethyl sulfate (0.91 mL, 9.6 mmol) dissolved in toluene (5 mL) was added slowly. The mixture was heated at 35°C for 20 hours. After cooling, the solids were removed by filtration and the solvent was concentrated *in vacuo*. The desired material was purified by chromatography on SiO₂ using 30% ethyl acetate / hexanes as eluent. MS (*m/z*): 248/250 (M+H)⁺; C₁₃H₁₃N₂OCl requires 249.

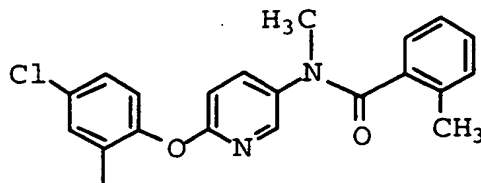
Example 15Preparation of 2,6-Dichloro-N-(6-(4-chloro-2-methylphenoxy)-pyridin-3-yl)-N-methylbenzamide

2-(4-Chloro-2-methyl-phenoxy)-5-(N-methylamino)pyridine (32 mg, 0.13 mmol) dissolved in methylene chloride (0.25 mL) was treated with triethylamine (0.026 mL, 0.18 mmol) followed by a solution of 2,6-dichlorobenzoyl chloride (31 mg, 0.15 mmol) dissolved in methylene chloride (0.15 mL). The reaction mixture was shaken at 22°C for 18 hours followed by quenching with saturated aqueous NH_4Cl and stirring for an additional 5 hours. The organic layer was separated and dried over Na_2SO_4 , then concentrated in vacuo. The crude product was purified by chromatography on SiO_2 using 1:1 ethyl acetate / hexane as eluent. MS (m/z): 422/424 ($\text{M}+\text{H}$)⁺; $\text{C}_{20}\text{H}_{15}\text{N}_2\text{O}_2\text{Cl}_3$, requires 422.

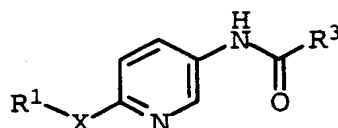
Example 16Preparation of 2-Chloro-N-(6-(4-chloro-2-methylphenoxy)-pyridin-3-yl)-N-methylbenzamide

2-Chloro-N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-N-methylbenzamide was prepared from 2-(4-Chloro-2-methyl-phenoxy)-5-(N-methylamino)pyridine and 2-chlorobenzoyl chloride in the same manner as 2,6-Dichloro-N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-N-methylbenzamide was prepared.

57

Example 17Preparation of 2-Methyl-N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-N-methyl-benzamide

- 5 2-Methyl-N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-N-methyl-benzamide was prepared from 2-(4-Chloro-2-methyl-phenoxy)-5-(N-methylamino)pyridine and 2-methylbenzoyl chloride in the same manner as 2,6-Dichloro-N-(6-(4-chloro-2-methyl-phenoxy)-pyridin-3-yl)-N-methyl-benzamide was prepared.
- 10

Example 18

15 General procedure for the synthesis of 2-substituted-5-acylamino-pyridines

- A solution of the 2-substituted-5-aminopyridine (10 mmol), triethylamine (20 mmol) and an acid chloride (20 mmol) in ethanol free chloroform (250 mL) was shaken for 16 hours. The mixture was then diluted with saturated aqueous sodium hydrogencarbonate (50 mL) and dichloromethane (500 mL), and shaken for 30 min. The mixture was then filtered through anhydrous magnesium sulfate, washing with dichloromethane (250 mL). Concentration of the filtrate under reduced pressure afforded the desired 2-substituted-5-acylamino-pyridines.
- 20
- 25

- The compounds listed in Table 6 were prepared from substituted 5-aminopyridine compounds and the appropriate acid chloride according to the general procedure above.
- 30

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Table 6

| <u>R¹X</u> | <u>R³</u> | <u>MS</u> <u>(m/z)</u> |
|--------------------------|---------------------------|---------------------------|
| 4-chloro-2-methylphenoxy | 4-biphenyl | 415 |
| 4-chloro-2-methylphenoxy | 3,4-dimethoxyphenyl | 319 |
| 4-chloro-2-methylphenoxy | 2-(trifluoromethyl)phenyl | 407 |
| 4-chloro-2-methylphenoxy | 2,4-difluorophenyl | 375 |
| 4-chloro-2-methylphenoxy | 4-cyanophenyl | 364 |
| 4-chloro-2-methylphenoxy | 3-(trifluoromethyl)phenyl | 407 |
| 4-chloro-2-methylphenoxy | 3-cyanophenyl | 364 |
| 4-chloro-2-methylphenoxy | 2-naphthyl | 389 |
| 4-chloro-2-methylphenoxy | 2-methoxyphenyl | 369 |
| 4-chloro-2-methylphenoxy | 3,4,5-trimethylphenyl | 429 |
| 4-chloro-2-methylphenoxy | 4-nitrophenyl | 384 |
| 4-chloro-2-methylphenoxy | 3,4-dichlorophenyl | 408 |
| 4-chloro-2-methylphenoxy | 5-nitrofuran-2-yl | 374 |
| 4-chloro-2-methylphenoxy | 3-bromophenyl | 418 |

| | | |
|--------------------------|------------------------------------|-----|
| 4-chloro-2-methylphenoxy | 3-pyridyl | 340 |
| 4-chloro-2-methylphenoxy | 2-ethoxynaphth-1-yl | 433 |
| 4-chloro-2-methylphenoxy | 2,3-dichlorophenyl | 408 |
| 4-chloro-2-methylphenoxy | 3-nitrophenyl | 384 |
| 4-chloro-2-methylphenoxy | 6-chloropyrid-3-yl | 374 |
| 4-chloro-2-methylphenoxy | 4-(trifluoromethoxy)phenyl | 423 |
| 4-chloro-2-methylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 425 |
| 4-chloro-2-methylphenoxy | 2-acetoxyphenyl | 397 |
| 4-chloro-2-methylphenoxy | 5-methylisoxazol-3-yl | 344 |
| 4-chloro-2-methylphenoxy | 2-(phenylthio)pyrid-3-yl | 448 |
| 4-chloro-2-methylphenoxy | 2-(trifluoromethoxy)phenyl | 423 |
| 4-chloro-2-methylphenoxy | 1-phenyl-5-propyl-pyrazin-4-yl | 447 |
| 4-chloro-2-methylphenoxy | 2-ethoxyphenyl | 383 |
| 4-chloro-2-methylphenoxy | 3-chlorothien-2-yl | 379 |
| 4-chloro-2-methylphenoxy | 3-bromothien-2-yl | 424 |

| | | |
|--------------------------|---|-----|
| 4-chloro-2-methylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 399 |
| 4-chloro-2-methylphenoxy | 3,5-dichlorophenyl | 408 |
| 4-chloro-2-methylphenoxy | 2-(propylthio)pyridin-3-yl | 414 |
| 4-chloro-2-methylphenoxy | 2-(ethylthio)pyridin-3-yl | 400 |
| 4-chloro-2-methylphenoxy | 3-bromopyridin-5-yl | 419 |
| 4-chloro-2-methylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 361 |
| 4-chloro-2-methylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 399 |
| 4-chloro-2-methylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 429 |
| 4-chloro-2-methylphenoxy | 4-chlorophenyl | 373 |
| 4-chloro-2-methylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 420 |
| 4-chloro-2-methylphenoxy | benzo[b]thiophen-2-yl | 395 |
| 4-chloro-2-methylphenoxy | 3,4-dimethylphenyl | 367 |
| 4-chloro-2-methylphenoxy | 2-(phenoxy)pyridin-3-yl | 432 |
| 4-chloro-2-methylphenoxy | 2-(methylthio)pyridin-3-yl | 386 |
| 4-chloro-2-methylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 420 |

| | | |
|--------------------------|--|-----|
| 4-chloro-2-methylphenoxy | 4-chloro-1,3-dimethyl pyrazolo[3,4-b]pyridin-3-yl | 442 |
| 4-chloro-2-methylphenoxy | 2-chloro-6-methylpyridin-4-yl | 388 |
| 4-chloro-2-methylphenoxy | 3,5-dimethylisoxazol-4-yl | 358 |
| 4-chloro-2-methylphenoxy | 1-naphthyl | 389 |
| 4-chloro-2-methylphenoxy | 2-fluorophenyl | 357 |
| 4-chloro-2-methylphenoxy | 4-propylphenyl | 381 |
| 4-chloro-2-methylphenoxy | 4-(trifluoromethyl)phenyl | 407 |
| 4-chloro-2-methylphenoxy | 3-fluorophenyl | 357 |
| 4-chloro-2-methylphenoxy | 2,6-difluorophenyl | 375 |
| 4-chloro-2-methylphenoxy | 2-chlorophenyl | 373 |
| 4-chloro-2-methylphenoxy | 3-(chloromethyl)phenyl | 387 |
| 4-chloro-2-methylphenoxy | 4-(2-(2-methyl) propyl)phenyl | 395 |
| 4-chloro-2-methylphenoxy | 3-chlorophenyl | 373 |
| 4-chloro-2-methylphenoxy | 2-nitrophenyl | 384 |
| 4-chloro-2-methylphenoxy | 3,5-dimethoxyphenyl | 399 |

| | | |
|--------------------------|------------------------------------|-----|
| 4-chloro-2-methylphenoxy | 2,6-dichlorophenyl | 408 |
| 4-chloro-2-methylphenoxy | 2,4-dichlorophenyl | 408 |
| 4-chloro-2-methylphenoxy | 4-fluorophenyl | 357 |
| 4-chloro-2-methylphenoxy | 4-butylphenyl | 395 |
| 4-chloro-2-methylphenoxy | 2-methylphenyl | 353 |
| 4-chloro-2-methylphenoxy | phenyl | 339 |
| 4-chloro-2-methylphenoxy | 4-ethylphenyl | 367 |
| 4-chloro-2-methylphenoxy | 2,3-difluorophenyl | 375 |
| 4-chloro-2-methylphenoxy | 2,6-dimethoxyphenyl | 399 |
| 4-chloro-2-methylphenoxy | 2,5-difluorophenyl | 375 |
| 4-chloro-2-methylphenoxy | 4-ethoxyphenyl | 383 |
| 4-chloro-2-methylphenoxy | 2,4,6-trichlorophenyl | 442 |
| 4-chloro-2-methylphenoxy | 3-methylphenyl | 353 |
| 4-chloro-2-methylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 425 |
| 4-chloro-2-methylphenoxy | 3-methoxyphenyl | 369 |
| 4-chloro-2-methylphenoxy | thien-2-yl | 345 |

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| 4-chloro-2-methylphenoxy | 2-bromophenyl | 418 |
| 4-chloro-2-methylphenoxy | 4-bromophenyl | 418 |
| 4-chloro-2-methylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 425 |
| 4-chloro-2-methylphenoxy | 3-(trifluoromethoxy)phenyl | 423 |
| 4-chloro-2-methylphenoxy | 9-fluorenon-4-yl | 441 |
| 4-chloro-2-methylphenoxy | isoxazol-5-yl | 330 |
| 4-chloro-2-methylphenoxy | benzofuroxan-5-yl | 397 |
| 4-chloro-2-methylphenoxy | 2-chloropyrid-3-yl | 374 |
| 4-chloro-2-methylphenoxy | 3,5-difluorophenyl | 375 |
| 4-chloro-2-methylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 446 |
| 4-chloro-2-methylphenoxy | pyridin-4-yl | 340 |
| 4-chloro-2-methylphenoxy | anthraquinon-2-yl | 469 |
| 4-chloro-2-methylphenoxy | 2-iodophenyl | 465 |
| 1-naphthoxy | 4-biphenyl | 416 |
| 1-naphthoxy | 3,4-dimethoxyphenyl | 400 |
| 1-naphthoxy | 2-(trifluoromethyl)phenyl | 408 |
| 1-naphthoxy | 2,4-difluorophenyl | 376 |
| 1-naphthoxy | 4-cyanophenyl | 365 |

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| 1-naphthoxy | 3-(trifluoromethyl)phenyl | 408 |
| 1-naphthoxy | 3-cyanophenyl | 365 |
| 1-naphthoxy | 2-naphthyl | 390 |
| 1-naphthoxy | 2-methoxyphenyl | 370 |
| 1-naphthoxy | 3,4,5-trimethylphenyl | 430 |
| 1-naphthoxy | 4-nitrophenyl | 385 |
| 1-naphthoxy | 3,4-dichlorophenyl | 409 |
| 1-naphthoxy | 5-nitrofuran-2-yl | 375 |
| 1-naphthoxy | 3-bromophenyl | 419 |
| 1-naphthoxy | 3-pyridyl | 341 |
| 1-naphthoxy | 2-ethoxynaphth-1-yl | 334 |
| 1-naphthoxy | 2,3-dichlorophenyl | 409 |
| 1-naphthoxy | 3-nitrophenyl | 385 |
| 1-naphthoxy | 6-chloropyrid-3-yl | 376 |
| 1-naphthoxy | 4-(trifluoromethoxy)phenyl | 424 |
| 1-naphthoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 426 |
| 1-naphthoxy | 3-bromothiophenyl | 425 |
| 1-naphthoxy | 2-acetoxyphenyl | 398 |
| 1-naphthoxy | 5-methylisoxazol-3-yl | 345 |
| 1-naphthoxy | 2-(phenylthio)pyrid-3-yl | 449 |
| 1-naphthoxy | 2-(trifluoromethoxy)phenyl | 424 |
| 1-naphthoxy | 1-phenyl-5-propylpyrazin-4-yl | 448 |
| 1-naphthoxy | 2-ethoxyphenyl | 384 |
| 1-naphthoxy | 3-chlorothien-2-yl | 381 |
| 1-naphthoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 400 |

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| 1-naphthoxy | 3,5-dichlorophenyl | 409 |
| 1-naphthoxy | 2-(propylthio)pyridin-3-yl | 415 |
| 1-naphthoxy | 2-(ethylthio)pyridin-3-yl | 401 |
| 1-naphthoxy | 3-bromopyridin-5-yl | 420 |
| 1-naphthoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 362 |
| 1-naphthoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 400 |
| 1-naphthoxy | 3-chlorobenzo[b]thiophen-2-yl | 431 |
| 1-naphthoxy | 4-chlorophenyl | 375 |
| 1-naphthoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 421 |
| 1-naphthoxy | benzo[b]thiophen-2-yl | 396 |
| 1-naphthoxy | 3,4-dimethylphenyl | 368 |
| 1-naphthoxy | 2-(phenoxy)pyridin-3-yl | 433 |
| 1-naphthoxy | 2-(methylthio)pyridin-3-yl | 387 |
| 1-naphthoxy | 5-methyl-3-phenylisoxazol-4-yl | 421 |
| 1-naphthoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 444 |
| 1-naphthoxy | 2-chloro-6-methylpyridin-4-yl | 390 |
| 1-naphthoxy | 3,5-dimethylisoxazol-4-yl | 359 |
| 1-naphthoxy | 1-naphthyl | 390 |
| 1-naphthoxy | 2-fluorophenyl | 358 |
| 1-naphthoxy | 4-propylphenyl | 382 |
| 1-naphthoxy | 4-(trifluoromethyl)phenyl | 408 |
| 1-naphthoxy | 3-fluorophenyl | 358 |

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| 1-naphthoxy | 2,6-difluorophenyl | 376 |
| 1-naphthoxy | 2-chlorophenyl | 375 |
| 1-naphthoxy | 3-(chloromethyl)phenyl | 389 |
| 1-naphthoxy | 4-(2-(2-methyl)propyl)phenyl | 396 |
| 1-naphthoxy | 3-chlorophenyl | 375 |
| 1-naphthoxy | 2-nitrophenyl | 385 |
| 1-naphthoxy | 3,5-dimethoxyphenyl | 400 |
| 1-naphthoxy | 2,6-dichlorophenyl | 409 |
| 1-naphthoxy | 2,4-dichlorophenyl | 409 |
| 1-naphthoxy | 4-fluorophenyl | 358 |
| 1-naphthoxy | 4-butylphenyl | 396 |
| 1-naphthoxy | 2-methylphenyl | 354 |
| 1-naphthoxy | phenyl | 340 |
| 1-naphthoxy | 4-ethylphenyl | 368 |
| 1-naphthoxy | 2,3-difluorophenyl | 376 |
| 1-naphthoxy | 2,6-dimethoxyphenyl | 400 |
| 1-naphthoxy | 3,4-difluorophenyl | 376 |
| 1-naphthoxy | 2,5-difluorophenyl | 376 |
| 1-naphthoxy | 4-ethoxyphenyl | 384 |
| 1-naphthoxy | 2,4,6-trichlorophenyl | 444 |
| 1-naphthoxy | 3-methylphenyl | 354 |
| 1-naphthoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 426 |
| 1-naphthoxy | 3-methoxyphenyl | 370 |
| 1-naphthoxy | thien-2-yl | 346 |
| 1-naphthoxy | 2-bromophenyl | 419 |
| 1-naphthoxy | 4-bromophenyl | 419 |

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| 1-naphthoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 426 |
| 1-naphthoxy | 3-(trifluoromethoxy)phenyl | 424 |
| 1-naphthoxy | 9-fluorenon-4-yl | 442 |
| 1-naphthoxy | isoxazol-5-yl | 331 |
| 1-naphthoxy | benzofuroxan-5-yl | 398 |
| 1-naphthoxy | 2-chloropyrid-3-yl | 376 |
| 1-naphthoxy | 3,5-difluorophenyl | 376 |
| 1-naphthoxy | 2-(4-methylphenoxy)pyridin-3-yl | 447 |
| 1-naphthoxy | pyridin-4-yl | 341 |
| 1-naphthoxy | anthraquinon-2-yl | 470 |
| 1-naphthoxy | 2-iodophenyl | 466 |
| 2-(2-propyl)phenoxy | 4-biphenyl | 408 |
| 2-(2-propyl)phenoxy | 3,4-dimethoxyphenyl | 392 |
| 2-(2-propyl)phenoxy | 2-(trifluoromethyl)phenyl | 400 |
| 2-(2-propyl)phenoxy | 2,4-difluorophenyl | 368 |
| 2-(2-propyl)phenoxy | 4-cyanophenyl | 357 |
| 2-(2-propyl)phenoxy | 3-(trifluoromethyl)phenyl | 400 |
| 2-(2-propyl)phenoxy | 3-cyanophenyl | 357 |
| 2-(2-propyl)phenoxy | 2-naphthyl | 382 |
| 2-(2-propyl)phenoxy | 2-methoxyphenyl | 362 |
| 2-(2-propyl)phenoxy | 3,4,5,-trimethylphenyl | 422 |
| 2-(2-propyl)phenoxy | 4-nitrophenyl | 377 |
| 2-(2-propyl)phenoxy | 3,4-dichlorophenyl | 401 |
| 2-(2-propyl)phenoxy | 5-nitrofuran-2-yl | 367 |
| 2-(2-propyl)phenoxy | 3-bromophenyl | 411 |
| 2-(2-propyl)phenoxy | 3-pyridyl | 333 |

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| 2-(2-propyl)phenoxy | 2-ethoxynaphth-1-yl | 426 |
| 2-(2-propyl)phenoxy | 2,3-dichlorophenyl | 401 |
| 2-(2-propyl)phenoxy | 3-nitrophenyl | 377 |
| 2-(2-propyl)phenoxy | 6-chloropyrid-3-yl | 368 |
| 2-(2-propyl)phenoxy | 4-(trifluoromethoxy)phenyl | 416 |
| 2-(2-propyl)phenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 418 |
| 2-(2-propyl)phenoxy | 3-bromothiophenyl | 417 |
| 2-(2-propyl)phenoxy | 2-acetoxyphenyl | 390 |
| 2-(2-propyl)phenoxy | 5-methylisoxazol-3-yl | 337 |
| 2-(2-propyl)phenoxy | 2-(phenylthio)pyrid-3-yl | 442 |
| 2-(2-propyl)phenoxy | 2-(trifluoromethoxy)phenyl | 416 |
| 2-(2-propyl)phenoxy | 1-phenyl-5-propylpyrazin-4-yl | 441 |
| 2-(2-propyl)phenoxy | 2-ethoxyphenyl | 376 |
| 2-(2-propyl)phenoxy | 3-chlorothien-2-yl | 373 |
| 2-(2-propyl)phenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 392 |
| 2-(2-propyl)phenoxy | 3,5-dichlorophenyl | 401 |
| 2-(2-propyl)phenoxy | 2-(propylthio)pyridin-3-yl | 407 |
| 2-(2-propyl)phenoxy | 2-(ethylthio)pyridin-3-yl | 393 |
| 2-(2-propyl)phenoxy | 3-bromopyridin-5-yl | 412 |
| 2-(2-propyl)phenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 354 |
| 2-(2-propyl)phenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 392 |
| 2-(2-propyl)phenoxy | 3-chlorobenzo[b]thiophen-2-yl | 423 |
| 2-(2-propyl)phenoxy | 4-chlorophenyl | 367 |

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| 2-(2-propyl)phenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 413 |
| 2-(2-propyl)phenoxy | benzo[b]thiophen-2-yl | 388 |
| 2-(2-propyl)phenoxy | 3,4-dimethylphenyl | 360 |
| 2-(2-propyl)phenoxy | 2-(phenoxy)pyridin-3-yl | 425 |
| 2-(2-propyl)phenoxy | 2-(methylthio)pyridin-3-yl | 379 |
| 2-(2-propyl)phenoxy | 5-methyl-3-phenylisoxazol-4-yl | 413 |
| 2-(2-propyl)phenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 436 |
| 2-(2-propyl)phenoxy | 2-chloro-6-methylpyridin-4-yl | 382 |
| 2-(2-propyl)phenoxy | 3,5-dimethylisoxazol-4-yl | 351 |
| 2-(2-propyl)phenoxy | 1-naphthyl | 382 |
| 2-(2-propyl)phenoxy | 2-fluorophenyl | 350 |
| 2-(2-propyl)phenoxy | 4-propylphenyl | 374 |
| 2-(2-propyl)phenoxy | 4-(trifluoromethyl)phenyl | 400 |
| 2-(2-propyl)phenoxy | 3-fluorophenyl | 350 |
| 2-(2-propyl)phenoxy | 2,6-difluorophenyl | 368 |
| 2-(2-propyl)phenoxy | 2-chlorophenyl | 367 |
| 2-(2-propyl)phenoxy | 3-(chloromethyl)phenyl | 381 |
| 2-(2-propyl)phenoxy | 4-(2-(2-methyl)propyl)phenyl | 388 |
| 2-(2-propyl)phenoxy | 3-chlorophenyl | 367 |
| 2-(2-propyl)phenoxy | 2-nitrophenyl | 377 |
| 2-(2-propyl)phenoxy | 3,5-dimethoxyphenyl | 392 |
| 2-(2-propyl)phenoxy | 2,6-dichlorophenyl | 401 |
| 2-(2-propyl)phenoxy | 2,4-dichlorophenyl | 401 |

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| 2-(2-propyl)phenoxy | 4-fluorophenyl | |
| 2-(2-propyl)phenoxy | 4-butylphenyl | |
| 2-(2-propyl)phenoxy | 2-methylphenyl | |
| 2-(2-propyl)phenoxy | phenyl | |
| 2-(2-propyl)phenoxy | 4-ethylphenyl | |
| 2-(2-propyl)phenoxy | 2,3-difluorophenyl | |
| 2-(2-propyl)phenoxy | 2,6-dimethoxyphenyl | 392 |
| 2-(2-propyl)phenoxy | 3,4-difluorophenyl | 368 |
| 2-(2-propyl)phenoxy | 2,5-difluorophenyl | 368 |
| 2-(2-propyl)phenoxy | 4-ethoxyphenyl | 376 |
| 2-(2-propyl)phenoxy | 2,4,6-trichlorophenyl | 436 |
| 2-(2-propyl)phenoxy | 3-methylphenyl | 346 |
| 2-(2-propyl)phenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 418 |
| 2-(2-propyl)phenoxy | 3-methoxyphenyl | 362 |
| 2-(2-propyl)phenoxy | thien-2-yl | 338 |
| 2-(2-propyl)phenoxy | 2-bromophenyl | 411 |
| 2-(2-propyl)phenoxy | 4-bromophenyl | 411 |
| 2-(2-propyl)phenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 418 |
| 2-(2-propyl)phenoxy | 3-(trifluoromethoxy)phenyl | 416 |
| 2-(2-propyl)phenoxy | 9-fluorenon-4-yl | 434 |
| 2-(2-propyl)phenoxy | isoxazol-5-yl | 323 |
| 2-(2-propyl)phenoxy | benzofuroxan-5-yl | 390 |
| 2-(2-propyl)phenoxy | 2-chloropyrid-3-yl | 368 |
| 2-(2-propyl)phenoxy | 3,5-difluorophenyl | 368 |
| 2-(2-propyl)phenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 439 |

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| 2-(2-propyl)phenoxy | pyridin-4-yl | 333 |
| 2-(2-propyl)phenoxy | anthraquinon-2-yl | 462 |
| 2-(2-propyl)phenoxy | 2-iodophenyl | 458 |
| 3-fluoro-5-methylphenoxy | 4-biphenyl | 398 |
| 3-fluoro-5-methylphenoxy | 3,4-dimethoxyphenyl | 382 |
| 3-fluoro-5-methylphenoxy | 2-(trifluoromethyl)phenyl | 390 |
| 3-fluoro-5-methylphenoxy | 2,4-difluorophenyl | 358 |
| 3-fluoro-5-methylphenoxy | 4-cyanophenyl | 347 |
| 3-fluoro-5-methylphenoxy | 3-(trifluoromethyl)phenyl | 390 |
| 3-fluoro-5-methylphenoxy | 3-cyanophenyl | 347 |
| 3-fluoro-5-methylphenoxy | 2-naphthyl | 372 |
| 3-fluoro-5-methylphenoxy | 2-methoxyphenyl | 352 |
| 3-fluoro-5-methylphenoxy | 3,4,5,-trimethylphenyl | 412 |
| 3-fluoro-5-methylphenoxy | 4-nitrophenyl | 367 |
| 3-fluoro-5-methylphenoxy | 3,4-dichlorophenyl | 391 |
| 3-fluoro-5-methylphenoxy | 5-nitrofuran-2-yl | 357 |
| 3-fluoro-5-methylphenoxy | 3-bromophenyl | 401 |

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| 3-fluoro-5-methylphenoxy | 3-pyridyl | 323 |
| 3-fluoro-5-methylphenoxy | 2-ethoxynaphth-1-yl | 416 |
| 3-fluoro-5-methylphenoxy | 2,3-dichlorophenyl | 391 |
| 3-fluoro-5-methylphenoxy | 3-nitrophenyl | 367 |
| 3-fluoro-5-methylphenoxy | 6-chloropyrid-3-yl | 358 |
| 3-fluoro-5-methylphenoxy | 4-(trifluoromethoxy)phenyl | 406 |
| 3-fluoro-5-methylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 408 |
| 3-fluoro-5-methylphenoxy | 3-bromothieryl | 407 |
| 3-fluoro-5-methylphenoxy | 2-acetoxyphenyl | 380 |
| 3-fluoro-5-methylphenoxy | 5-methylisoxazol-3-yl | 327 |
| 3-fluoro-5-methylphenoxy | 2-(phenylthio)pyrid-3-yl | 431 |
| 3-fluoro-5-methylphenoxy | 2-(trifluoromethoxy)phenyl | 406 |
| 3-fluoro-5-methylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 430 |
| 3-fluoro-5-methylphenoxy | 2-ethoxyphenyl | 366 |
| 3-fluoro-5-methylphenoxy | 3-chlorothien-2-yl | 363 |
| 3-fluoro-5-methylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 382 |

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| 3-fluoro-5-methylphenoxy | 3,5-dichlorophenyl | 391 |
| 3-fluoro-5-methylphenoxy | 2-(propylthio)pyridin-3-yl | 397 |
| 3-fluoro-5-methylphenoxy | 2-(ethylthio)pyridin-3-yl | 383 |
| 3-fluoro-5-methylphenoxy | 3-bromopyridin-5-yl | 402 |
| 3-fluoro-5-methylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 344 |
| 3-fluoro-5-methylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 382 |
| 3-fluoro-5-methylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 413 |
| 3-fluoro-5-methylphenoxy | 4-chlorophenyl | 357 |
| 3-fluoro-5-methylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 403 |
| 3-fluoro-5-methylphenoxy | benzo[b]thiophen-2-yl | 378 |
| 3-fluoro-5-methylphenoxy | 3,4-dimethylphenyl | 350 |
| 3-fluoro-5-methylphenoxy | 2-(phenoxy)pyridin-3-yl | 415 |
| 3-fluoro-5-methylphenoxy | 2-(methylthio)pyridin-3-yl | 369 |
| 3-fluoro-5-methylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 403 |
| 3-fluoro-5-methylphenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 426 |

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| 3-fluoro-5-methylphenoxy | 2-chloro-6-methylpyridin-4-yl | 372 |
| 3-fluoro-5-methylphenoxy | 3,5-dimethylisoxazol-4-yl | 341 |
| 3-fluoro-5-methylphenoxy | 1-naphthyl | 372 |
| 3-fluoro-5-methylphenoxy | 2-fluorophenyl | 340 |
| 3-fluoro-5-methylphenoxy | 4-propylphenyl | 364 |
| 3-fluoro-5-methylphenoxy | 4-(trifluoromethyl)phenyl | 390 |
| 3-fluoro-5-methylphenoxy | 3-fluorophenyl | 340 |
| 3-fluoro-5-methylphenoxy | 2,6-difluorophenyl | 358 |
| 3-fluoro-5-methylphenoxy | 2-chlorophenyl | 357 |
| 3-fluoro-5-methylphenoxy | 3-(chloromethyl)phenyl | 371 |
| 3-fluoro-5-methylphenoxy | 4-(2-(2-methyl)propyl)phenyl | 378 |
| 3-fluoro-5-methylphenoxy | 3-chlorophenyl | 357 |
| 3-fluoro-5-methylphenoxy | 2-nitrophenyl | 367 |
| 3-fluoro-5-methylphenoxy | 3,5-dimethoxyphenyl | 382 |
| 3-fluoro-5-methylphenoxy | 2,6-dichlorophenyl | 391 |

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| 3-fluoro-5-methylphenoxy | 2,4-dichlorophenyl | 391 |
| 3-fluoro-5-methylphenoxy | 4-fluorophenyl | 340 |
| 3-fluoro-5-methylphenoxy | 4-butylphenyl | 378 |
| 3-fluoro-5-methylphenoxy | 2-methylphenyl | 336 |
| 3-fluoro-5-methylphenoxy | phenyl | 322 |
| 3-fluoro-5-methylphenoxy | 4-ethylphenyl | 350 |
| 3-fluoro-5-methylphenoxy | 2,3-difluorophenyl | 358 |
| 3-fluoro-5-methylphenoxy | 2,6-dimethoxyphenyl | 382 |
| 3-fluoro-5-methylphenoxy | 3,4-difluorophenyl | 358 |
| 3-fluoro-5-methylphenoxy | 2,5-difluorophenyl | 358 |
| 3-fluoro-5-methylphenoxy | 4-ethoxyphenyl | 366 |
| 3-fluoro-5-methylphenoxy | 2,4,6-trichlorophenyl | 426 |
| 3-fluoro-5-methylphenoxy | 3-methylphenyl | 336 |
| 3-fluoro-5-methylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 408 |
| 3-fluoro-5-methylphenoxy | 3-methoxyphenyl | 352 |

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| 3-fluoro-5-methylphenoxy | thien-2-yl | 328 |
| 3-fluoro-5-methylphenoxy | 2-bromophenyl | 401 |
| 3-fluoro-5-methylphenoxy | 4-bromophenyl | 401 |
| 3-fluoro-5-methylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 408 |
| 3-fluoro-5-methylphenoxy | 3-(trifluoromethoxy)phenyl | 406 |
| 3-fluoro-5-methylphenoxy | 9-fluorenon-4-yl | 424 |
| 3-fluoro-5-methylphenoxy | isoxazol-5-yl | 313 |
| 3-fluoro-5-methylphenoxy | benzofuroxan-5-yl | 380 |
| 3-fluoro-5-methylphenoxy | 2-chloropyrid-3-yl | 358 |
| 3-fluoro-5-methylphenoxy | 3,5-difluorophenyl | 358 |
| 3-fluoro-5-methylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 429 |
| 3-fluoro-5-methylphenoxy | pyridin-4-yl | 323 |
| 3-fluoro-5-methylphenoxy | anthraquinon-2-yl | 452 |
| 3-fluoro-5-methylphenoxy | 2-iodophenyl | 448 |
| 2-methylpyrid-3-yloxy | 4-biphenyl | 381 |
| 2-methylpyrid-3-yloxy | 3,4-dimethoxyphenyl | 365 |

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| 2-methylpyrid-3-yloxy | 2-(trifluoromethyl)phenyl | 373 |
| 2-methylpyrid-3-yloxy | 2,4-difluorophenyl | 341 |
| 2-methylpyrid-3-yloxy | 4-cyanophenyl | 330 |
| 2-methylpyrid-3-yloxy | 3-(trifluoromethyl)phenyl | 373 |
| 2-methylpyrid-3-yloxy | 3-cyanophenyl | 330 |
| 2-methylpyrid-3-yloxy | 2-naphthyl | 355 |
| 2-methylpyrid-3-yloxy | 2-methoxyphenyl | 335 |
| 2-methylpyrid-3-yloxy | 3,4,5-trimethylphenyl | 395 |
| 2-methylpyrid-3-yloxy | 4-nitrophenyl | 350 |
| 2-methylpyrid-3-yloxy | 3,4-dichlorophenyl | 374 |
| 2-methylpyrid-3-yloxy | 5-nitrofuran-2-yl | 340 |
| 2-methylpyrid-3-yloxy | 3-bromophenyl | 384 |
| 2-methylpyrid-3-yloxy | 3-pyridyl | 306 |
| 2-methylpyrid-3-yloxy | 2-ethoxynaphth-1-yl | 399 |
| 2-methylpyrid-3-yloxy | 2,3-dichlorophenyl | 374 |
| 2-methylpyrid-3-yloxy | 3-nitrophenyl | 350 |
| 2-methylpyrid-3-yloxy | 6-chloropyrid-3-yl | 341 |
| 2-methylpyrid-3-yloxy | 4-(trifluoromethoxy)phenyl | 389 |
| 2-methylpyrid-3-yloxy | 2-fluoro-4-(trifluoromethyl)phenyl | 391 |
| 2-methylpyrid-3-yloxy | 3-bromothieryl | 390 |
| 2-methylpyrid-3-yloxy | 2-acetoxyphenyl | 363 |
| 2-methylpyrid-3-yloxy | 5-methylisoxazol-3-yl | 310 |
| 2-methylpyrid-3-yloxy | 2-(phenylthio)pyrid-3-yl | 414 |
| 2-methylpyrid-3-yloxy | 2-(trifluoromethoxy)phenyl | 389 |
| 2-methylpyrid-3-yloxy | 1-phenyl-5-propylpyrazin-4-yl | 413 |
| 2-methylpyrid-3-yloxy | 2-ethoxyphenyl | 349 |

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| 2-methylpyrid-3-yloxy | 3-chlorothien-2-yl | 346 |
| 2-methylpyrid-3-yloxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 365 |
| 2-methylpyrid-3-yloxy | 3,5-dichlorophenyl | 374 |
| 2-methylpyrid-3-yloxy | 2-(propylthio)pyridin-3-yl | 380 |
| 2-methylpyrid-3-yloxy | 2-(ethylthio)pyridin-3-yl | 366 |
| 2-methylpyrid-3-yloxy | 3-bromopyridin-5-yl | 385 |
| 2-methylpyrid-3-yloxy | 4-methyl-1,2,3-thiadiazol-5-yl | 327 |
| 2-methylpyrid-3-yloxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 365 |
| 2-methylpyrid-3-yloxy | 3-chlorobenzo[b]thiophen-2-yl | 396 |
| 2-methylpyrid-3-yloxy | 4-chlorophenyl | 340 |
| 2-methylpyrid-3-yloxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 386 |
| 2-methylpyrid-3-yloxy | benzo[b]thiophen-2-yl | 361 |
| 2-methylpyrid-3-yloxy | 3,4-dimethylphenyl | 333 |
| 2-methylpyrid-3-yloxy | 2-(phenoxy)pyridin-3-yl | 398 |
| 2-methylpyrid-3-yloxy | 2-(methylthio)pyridin-3-yl | 352 |
| 2-methylpyrid-3-yloxy | 5-methyl-3-phenylisoxazol-4-yl | 386 |
| 2-methylpyrid-3-yloxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 409 |
| 2-methylpyrid-3-yloxy | 2-chloro-6-methylpyridin-4-yl | 355 |
| 2-methylpyrid-3-yloxy | 3,5-dimethylisoxazol-4-yl | 324 |
| 2-methylpyrid-3-yloxy | 1-naphthyl | 355 |
| 2-methylpyrid-3-yloxy | 2-fluorophenyl | 323 |

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| 2-methylpyrid-3-yloxy | 4-propylphenyl | 347 |
| 2-methylpyrid-3-yloxy | 4-(trifluoromethyl)phenyl | 373 |
| 2-methylpyrid-3-yloxy | 3-fluorophenyl | 323 |
| 2-methylpyrid-3-yloxy | 2,6-difluorophenyl | 341 |
| 2-methylpyrid-3-yloxy | 2-chlorophenyl | 340 |
| 2-methylpyrid-3-yloxy | 3-(chloromethyl)phenyl | 354 |
| 2-methylpyrid-3-yloxy | 4-(2-(2-methyl)propyl)phenyl | 361 |
| 2-methylpyrid-3-yloxy | 3-chlorophenyl | 340 |
| 2-methylpyrid-3-yloxy | 2-nitrophenyl | 350 |
| 2-methylpyrid-3-yloxy | 3,5-dimethoxyphenyl | 365 |
| 2-methylpyrid-3-yloxy | 2,6-dichlorophenyl | 374 |
| 2-methylpyrid-3-yloxy | 2,4-dichlorophenyl | 374 |
| 2-methylpyrid-3-yloxy | 4-fluorophenyl | 323 |
| 2-methylpyrid-3-yloxy | 4-butylphenyl | 361 |
| 2-methylpyrid-3-yloxy | 2-methylphenyl | 319 |
| 2-methylpyrid-3-yloxy | phenyl | 305 |
| 2-methylpyrid-3-yloxy | 4-ethylphenyl | 333 |
| 2-methylpyrid-3-yloxy | 2,3-difluorophenyl | 341 |
| 2-methylpyrid-3-yloxy | 2,6-dimethoxyphenyl | 365 |
| 2-methylpyrid-3-yloxy | 3,4-difluorophenyl | 341 |
| 2-methylpyrid-3-yloxy | 2,5-difluorophenyl | 341 |
| 2-methylpyrid-3-yloxy | 4-ethoxyphenyl | 349 |
| 2-methylpyrid-3-yloxy | 2,4,6-trichlorophenyl | 409 |
| 2-methylpyrid-3-yloxy | 3-methylphenyl | 319 |
| 2-methylpyrid-3-yloxy | 2-fluoro-5-(trifluoromethyl)phenyl | 391 |
| 2-methylpyrid-3-yloxy | 3-methoxyphenyl | 335 |

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| 2-methylpyrid-3-yloxy | 2-bromophenyl | 384 |
| 2-methylpyrid-3-yloxy | 4-bromophenyl | 384 |
| 2-methylpyrid-3-yloxy | 4-fluoro-3-(trifluoromethyl)phenyl | 391 |
| 2-methylpyrid-3-yloxy | 3-(trifluoromethoxy)phenyl | 389 |
| 2-methylpyrid-3-yloxy | 9-fluorenon-4-yl | 407 |
| 2-methylpyrid-3-yloxy | isoxazol-5-yl | 296 |
| 2-methylpyrid-3-yloxy | benzofuroxan-5-yl | 363 |
| 2-methylpyrid-3-yloxy | 2-chloropyrid-3-yl | 341 |
| 2-methylpyrid-3-yloxy | 3,5-difluorophenyl | 341 |
| 2-methylpyrid-3-yloxy | 2-(4-methylphenoxy)pyridin-3-yl | 412 |
| 2-methylpyrid-3-yloxy | pyridin-4-yl | 306 |
| 2-methylpyrid-3-yloxy | anthraquinon-2-yl | 435 |
| 2-methylpyrid-3-yloxy | 2-iodophenyl | 431 |
| 4-chloro-2,5-dimethylphenoxy | 3,4-dimethoxyphenyl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 2-(trifluoromethyl)phenyl | 421 |
| 4-chloro-2,5-dimethylphenoxy | 2,4-difluorophenyl | 389 |
| 4-chloro-2,5-dimethylphenoxy | 3-(trifluoromethyl)phenyl | 421 |
| 4-chloro-2,5-dimethylphenoxy | 2-naphthyl | 403 |
| 4-chloro-2,5-dimethylphenoxy | 2-methoxyphenyl | 484 |
| 4-chloro-2,5-dimethylphenoxy | 3,4,5-trimethylphenyl | 443 |

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| 4-chloro-2,5-dimethylphenoxy | 3,4-dichlorophenyl | 422 |
| 4-chloro-2,5-dimethylphenoxy | 3-bromophenyl | 432 |
| 4-chloro-2,5-dimethylphenoxy | 3-pyridyl | 354 |
| 4-chloro-2,5-dimethylphenoxy | 2-ethoxynaphth-1-yl | 447 |
| 4-chloro-2,5-dimethylphenoxy | 2,3-dichlorophenyl | 422 |
| 4-chloro-2,5-dimethylphenoxy | 6-chloropyrid-3-yl | 388 |
| 4-chloro-2,5-dimethylphenoxy | 4-(trifluoromethoxy)phenyl | 437 |
| 4-chloro-2,5-dimethylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 439 |
| 4-chloro-2,5-dimethylphenoxy | 3-bromothieryl | 438 |
| 4-chloro-2,5-dimethylphenoxy | 2-acetoxyphenyl | 411 |
| 4-chloro-2,5-dimethylphenoxy | 5-methylisoxazol-3-yl | 358 |
| 4-chloro-2,5-dimethylphenoxy | 2-(phenylthio)pyrid-3-yl | 462 |
| 4-chloro-2,5-dimethylphenoxy | 2-(trifluoromethoxy)phenyl | 437 |
| 4-chloro-2,5-dimethylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 461 |
| 4-chloro-2,5-dimethylphenoxy | 2-ethoxyphenyl | 397 |
| 4-chloro-2,5-dimethylphenoxy | 3-chlorothieryl | 393 |

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| 4-chloro-2,5-dimethylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 3,5-dichlorophenyl | 422 |
| 4-chloro-2,5-dimethylphenoxy | 2-(propylthio)pyridin-3-yl | 428 |
| 4-chloro-2,5-dimethylphenoxy | 2-(ethylthio)pyridin-3-yl | 414 |
| 4-chloro-2,5-dimethylphenoxy | 3-bromopyridin-5-yl | 433 |
| 4-chloro-2,5-dimethylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 375 |
| 4-chloro-2,5-dimethylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 443 |
| 4-chloro-2,5-dimethylphenoxy | 4-chlorophenyl | 387 |
| 4-chloro-2,5-dimethylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 434 |
| 4-chloro-2,5-dimethylphenoxy | benzo[b]thiophen-2-yl | 409 |
| 4-chloro-2,5-dimethylphenoxy | 3,4-dimethylphenyl | 381 |
| 4-chloro-2,5-dimethylphenoxy | 2-(phenoxy)pyridin-3-yl | 446 |
| 4-chloro-2,5-dimethylphenoxy | 2-(methylthio)pyridin-3-yl | 400 |
| 4-chloro-2,5-dimethylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 434 |

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| 4-chloro-2,5-dimethylphenoxy | 4-chloro-1,3-dimethyl pyrazolo[3,4-b]pyridin-3-yl | 456 |
| 4-chloro-2,5-dimethylphenoxy | 2-chloro-6-methylpyridin-4-yl | 402 |
| 4-chloro-2,5-dimethylphenoxy | 3,5-dimethylisoxazol-4-yl | 372 |
| 4-chloro-2,5-dimethylphenoxy | 1-naphthyl | 403 |
| 4-chloro-2,5-dimethylphenoxy | 2-fluorophenyl | 371 |
| 4-chloro-2,5-dimethylphenoxy | 4-propylphenyl | 395 |
| 4-chloro-2,5-dimethylphenoxy | 3-fluorophenyl | 371 |
| 4-chloro-2,5-dimethylphenoxy | 2,6-difluorophenyl | 389 |
| 4-chloro-2,5-dimethylphenoxy | 2-chlorophenyl | 387 |
| 4-chloro-2,5-dimethylphenoxy | 3-(chloromethyl)phenyl | 401 |
| 4-chloro-2,5-dimethylphenoxy | 4-(2-(2-methyl)propyl)phenyl | 409 |
| 4-chloro-2,5-dimethylphenoxy | 3-chlorophenyl | 387 |
| 4-chloro-2,5-dimethylphenoxy | 3,5-dimethoxyphenyl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 2,6-dichlorophenyl | 422 |
| 4-chloro-2,5-dimethylphenoxy | 2,4-dichlorophenyl | 422 |

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| 4-chloro-2,5-dimethylphenoxy | 4-fluorophenyl | 371 |
| 4-chloro-2,5-dimethylphenoxy | 4-butylphenyl | 409 |
| 4-chloro-2,5-dimethylphenoxy | 2-methylphenyl | 367 |
| 4-chloro-2,5-dimethylphenoxy | phenyl | 353 |
| 4-chloro-2,5-dimethylphenoxy | 4-ethylphenyl | 381 |
| 4-chloro-2,5-dimethylphenoxy | 2,3-difluorophenyl | 389 |
| 4-chloro-2,5-dimethylphenoxy | 2,6-dimethoxyphenyl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 3,4-difluorophenyl | 389 |
| 4-chloro-2,5-dimethylphenoxy | 2,5-difluorophenyl | 389 |
| 4-chloro-2,5-dimethylphenoxy | 4-ethoxyphenyl | 397 |
| 4-chloro-2,5-dimethylphenoxy | 2,4,6-trichlorophenyl | 456 |
| 4-chloro-2,5-dimethylphenoxy | 3-methylphenyl | 367 |
| 4-chloro-2,5-dimethylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 439 |
| 4-chloro-2,5-dimethylphenoxy | 3-methoxyphenyl | 383 |
| 4-chloro-2,5-dimethylphenoxy | 2-bromophenyl | 432 |
| 4-chloro-2,5-dimethylphenoxy | 4-bromophenyl | 432 |

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| 4-chloro-2,5-dimethylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 439 |
| 4-chloro-2,5-dimethylphenoxy | 3-(trifluoromethoxy)phenyl | 437 |
| 4-chloro-2,5-dimethylphenoxy | 9-fluorenon-4-yl | 455 |
| 4-chloro-2,5-dimethylphenoxy | isoxazol-5-yl | 344 |
| 4-chloro-2,5-dimethylphenoxy | benzofuroxan-5-yl | 411 |
| 4-chloro-2,5-dimethylphenoxy | 2-chloropyrid-3-yl | 388 |
| 4-chloro-2,5-dimethylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 460 |
| 4-chloro-2,5-dimethylphenoxy | pyridin-4-yl | 354 |
| 4-chloro-2,5-dimethylphenoxy | anthraquinon-2-yl | 483 |
| 4-chloro-2,5-dimethylphenoxy | 2-iodophenyl | 479 |
| 4-chloro-2,5-dimethylphenoxy | 4-pentylphenyl | 423 |
| 4-chloro-2,5-dimethylphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 496 |
| 4-chloro-2,5-dimethylphenoxy | 2,6-dimethylphenyl | 381 |
| 4-chloro-2,5-dimethylphenoxy | 2,5-dimethoxyphenyl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 2,5-dichloropyridin-3-yl | 423 |
| 4-chloro-2,5-dimethylphenoxy | 2-chloro-6-methoxypyridin-4-yl | 418 |

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| 4-chloro-2,5-dimethylphenoxy | 2,3-dichloropyridin-5-yl | 423 |
| 4-chloro-2,5-dimethylphenoxy | 1-naphthyl | 417 |
| 4-chloro-2,5-dimethylphenoxy | 2,4-dimethoxyphenyl | 413 |
| 4-chloro-2,5-dimethylphenoxy | 3,5-bis(trifluoromethyl)phenyl | 489 |
| 4-chloro-2,5-dimethylphenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 480 |
| 4-chloro-2,5-dimethylphenoxy | pentafluorophenyl | 443 |
| 4-methoxyphenoxy | 3,4-dimethoxyphenyl | 380 |
| 4-methoxyphenoxy | 2-(trifluoromethyl)phenyl | 388 |
| 4-methoxyphenoxy | 2,4-difluorophenyl | 356 |
| 4-methoxyphenoxy | 3-(trifluoromethyl)phenyl | 388 |
| 4-methoxyphenoxy | 2-naphthyl | 370 |
| 4-methoxyphenoxy | 2-methoxyphenyl | 350 |
| 4-methoxyphenoxy | 3,4,5-trimethylphenyl | 410 |
| 4-methoxyphenoxy | 3,4-dichlorophenyl | 389 |
| 4-methoxyphenoxy | 3-bromophenyl | 399 |
| 4-methoxyphenoxy | 3-pyridyl | 321 |
| 4-methoxyphenoxy | 2-ethoxynaphth-1-yl | 414 |
| 4-methoxyphenoxy | 2,3-dichlorophenyl | 389 |
| 4-methoxyphenoxy | 6-chloropyrid-3-yl | 356 |
| 4-methoxyphenoxy | 4-(trifluoromethoxy)phenyl | 404 |
| 4-methoxyphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 406 |
| 4-methoxyphenoxy | 3-bromothieryl | 405 |

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| 4-methoxyphenoxy | 2-acetoxyphenyl | 378 |
| 4-methoxyphenoxy | 5-methylisoxazol-3-yl | 325 |
| 4-methoxyphenoxy | 2-(phenylthio)pyrid-3-yl | 429 |
| 4-methoxyphenoxy | 2-(trifluoromethoxy)phenyl | 404 |
| 4-methoxyphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 428 |
| 4-methoxyphenoxy | 2-ethoxyphenyl | 364 |
| 4-methoxyphenoxy | 3-chlorothien-2-yl | 361 |
| 4-methoxyphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 380 |
| 4-methoxyphenoxy | 3,5-dichlorophenyl | 389 |
| 4-methoxyphenoxy | 2-(propylthio)pyridin-3-yl | 395 |
| 4-methoxyphenoxy | 2-(ethylthio)pyridin-3-yl | 381 |
| 4-methoxyphenoxy | 3-bromopyridin-5-yl | 400 |
| 4-methoxyphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 342 |
| 4-methoxyphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 380 |
| 4-methoxyphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 411 |
| 4-methoxyphenoxy | 4-chlorophenyl | 355 |
| 4-methoxyphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 401 |
| 4-methoxyphenoxy | benzo[b]thiophen-2-yl | 376 |
| 4-methoxyphenoxy | 3,4-dimethylphenyl | 348 |
| 4-methoxyphenoxy | 2-(phenoxy)pyridin-3-yl | 413 |
| 4-methoxyphenoxy | 2-(methylthio)pyridin-3-yl | 367 |
| 4-methoxyphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 401 |

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| 4-methoxyphenoxy | 4-chloro-1,3-dimethyl pyrazolo[3,4-b]pyridin-3- yl | 424 |
| 4-methoxyphenoxy | 2-chloro-6-methylpyridin- 4-yl | 370 |
| 4-methoxyphenoxy | 3,5-dimethylisoxazol-4-yl | 339 |
| 4-methoxyphenoxy | 1-naphthyl | 370 |
| 4-methoxyphenoxy | 2-fluorophenyl | 338 |
| 4-methoxyphenoxy | 4-propylphenyl | 362 |
| 4-methoxyphenoxy | 3-fluorophenyl | 338 |
| 4-methoxyphenoxy | 2,6-difluorophenyl | 356 |
| 4-methoxyphenoxy | 2-chlorophenyl | 355 |
| 4-methoxyphenoxy | 3-(chloromethyl)phenyl | 369 |
| 4-methoxyphenoxy | 4-(2-(2- methyl)propyl)phenyl | 376 |
| 4-methoxyphenoxy | 3-chlorophenyl | 355 |
| 4-methoxyphenoxy | 3,5-dimethoxyphenyl | 380 |
| 4-methoxyphenoxy | 2,6-dichlorophenyl | 389 |
| 4-methoxyphenoxy | 2,4-dichlorophenyl | 389 |
| 4-methoxyphenoxy | 4-fluorophenyl | 338 |
| 4-methoxyphenoxy | 4-butylphenyl | 376 |
| 4-methoxyphenoxy | 2-methylphenyl | 334 |
| 4-methoxyphenoxy | phenyl | 320 |
| 4-methoxyphenoxy | 4-ethylphenyl | 348 |
| 4-methoxyphenoxy | 2,3-difluorophenyl | 356 |
| 4-methoxyphenoxy | 2,6-dimethoxyphenyl | 380 |
| 4-methoxyphenoxy | 3,4-difluorophenyl | 356 |
| 4-methoxyphenoxy | 2,5-difluorophenyl | 356 |
| 4-methoxyphenoxy | 4-ethoxyphenyl | 364 |

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| 4-methoxyphenoxy | 2,4,6-trichlorophenyl | 424 |
| 4-methoxyphenoxy | 3-methylphenyl | 334 |
| 4-methoxyphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 406 |
| 4-methoxyphenoxy | 3-methoxyphenyl | 350 |
| 4-methoxyphenoxy | 2-bromophenyl | 399 |
| 4-methoxyphenoxy | 4-bromophenyl | 399 |
| 4-methoxyphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 406 |
| 4-methoxyphenoxy | 3-(trifluoromethoxy)phenyl | 404 |
| 4-methoxyphenoxy | 9-fluorenon-4-yl | 422 |
| 4-methoxyphenoxy | isoxazol-5-yl | 311 |
| 4-methoxyphenoxy | benzofuroxan-5-yl | 378 |
| 4-methoxyphenoxy | 2-chloropyrid-3-yl | 356 |
| 4-methoxyphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 427 |
| 4-methoxyphenoxy | pyridin-4-yl | 321 |
| 4-methoxyphenoxy | anthraquinon-2-yl | 450 |
| 4-methoxyphenoxy | 2-iodophenyl | 446 |
| 4-methoxyphenoxy | 4-pentylphenyl | 390 |
| 4-methoxyphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 464 |
| 4-methoxyphenoxy | 2,6-dimethylphenyl | 348 |
| 4-methoxyphenoxy | 2,5-dimethoxyphenyl | 380 |
| 4-methoxyphenoxy | 2,5-dichloropyridin-3-yl | 390 |
| 4-methoxyphenoxy | 2-chloro-6-methoxypyridin-4-yl | 386 |
| 4-methoxyphenoxy | 2,3-dichloropyridin-5-yl | 390 |
| 4-methoxyphenoxy | 1-naphthyl | 384 |

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| 4-methoxyphenoxy | 2,4-dimethoxyphenyl | 380 |
| 4-methoxyphenoxy | 3,5-bis(trifluoromethyl) phenyl | 456 |
| 4-methoxyphenoxy | 2-(4- chlorophenoxy)pyridin-3-yl | 448 |
| 4-methoxyphenoxy | pentafluorophenyl | 410 |
| 2-(2-propoxy)phenoxy | 3,4-dimethoxyphenyl | 408 |
| 2-(2-propoxy)phenoxy | 2-(trifluoromethyl)phenyl | 416 |
| 2-(2-propoxy)phenoxy | 2,4-difluorophenyl | 384 |
| 2-(2-propoxy)phenoxy | 3-(trifluoromethyl)phenyl | 416 |
| 2-(2-propoxy)phenoxy | 2-naphthyl | 398 |
| 2-(2-propoxy)phenoxy | 2-methoxyphenyl | 378 |
| 2-(2-propoxy)phenoxy | 3,4,5-trimethylphenyl | 438 |
| 2-(2-propoxy)phenoxy | 3,4-dichlorophenyl | 417 |
| 2-(2-propoxy)phenoxy | 3-bromophenyl | 427 |
| 2-(2-propoxy)phenoxy | 3-pyridyl | 349 |
| 2-(2-propoxy)phenoxy | 2-ethoxynaphth-1-yl | 442 |
| 2-(2-propoxy)phenoxy | 2,3-dichlorophenyl | 417 |
| 2-(2-propoxy)phenoxy | 6-chloropyrid-3-yl | 384 |
| 2-(2-propoxy)phenoxy | 4-(trifluoromethoxy)phenyl | 432 |
| 2-(2-propoxy)phenoxy | 2-fluoro-4- (trifluoromethyl)phenyl | 434 |
| 2-(2-propoxy)phenoxy | 3-bromothieryl | 433 |
| 2-(2-propoxy)phenoxy | 2-acetoxyphenyl | 406 |
| 2-(2-propoxy)phenoxy | 5-methylisoxazol-3-yl | 353 |
| 2-(2-propoxy)phenoxy | 2-(phenylthio)pyrid-3-yl | 458 |
| 2-(2-propoxy)phenoxy | 2-(trifluoromethoxy)phenyl | 432 |
| 2-(2-propoxy)phenoxy | 1-phenyl-5-propylpyrazin- 4-yl | 457 |

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| 2-(2-propoxy)phenoxy | 2-ethoxyphenyl | 392 |
| 2-(2-propoxy)phenoxy | 3-chlorothien-2-yl | 389 |
| 2-(2-propoxy)phenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 408 |
| 2-(2-propoxy)phenoxy | 3,5-dichlorophenyl | 417 |
| 2-(2-propoxy)phenoxy | 2-(propylthio)pyridin-3-yl | 423 |
| 2-(2-propoxy)phenoxy | 2-(ethylthio)pyridin-3-yl | 409 |
| 2-(2-propoxy)phenoxy | 3-bromopyridin-5-yl | 428 |
| 2-(2-propoxy)phenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 370 |
| 2-(2-propoxy)phenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 408 |
| 2-(2-propoxy)phenoxy | 3-chlorobenzo[b]thiophen-2-yl | 439 |
| 2-(2-propoxy)phenoxy | 4-chlorophenyl | 383 |
| 2-(2-propoxy)phenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 429 |
| 2-(2-propoxy)phenoxy | benzo[b]thiophen-2-yl | 404 |
| 2-(2-propoxy)phenoxy | 3,4-dimethylphenyl | 376 |
| 2-(2-propoxy)phenoxy | 2-(phenoxy)pyridin-3-yl | 441 |
| 2-(2-propoxy)phenoxy | 2-(methylthio)pyridin-3-yl | 395 |
| 2-(2-propoxy)phenoxy | 5-methyl-3-phenylisoxazol-4-yl | 429 |
| 2-(2-propoxy)phenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 452 |
| 2-(2-propoxy)phenoxy | 2-chloro-6-methylpyridin-4-yl | 398 |
| 2-(2-propoxy)phenoxy | 3,5-dimethylisoxazol-4-yl | 367 |
| 2-(2-propoxy)phenoxy | 1-naphthyl | 398 |

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| 2-(2-propoxy)phenoxy | 2-fluorophenyl | 366 |
| 2-(2-propoxy)phenoxy | 4-propylphenyl | 390 |
| 2-(2-propoxy)phenoxy | 3-fluorophenyl | 366 |
| 2-(2-propoxy)phenoxy | 2,6-difluorophenyl | 384 |
| 2-(2-propoxy)phenoxy | 2-chlorophenyl | 383 |
| 2-(2-propoxy)phenoxy | 3-(chloromethyl)phenyl | 397 |
| 2-(2-propoxy)phenoxy | 4-(2-(2-methyl)propyl)phenyl | 404 |
| 2-(2-propoxy)phenoxy | 3-chlorophenyl | 383 |
| 2-(2-propoxy)phenoxy | 3,5-dimethoxyphenyl | 408 |
| 2-(2-propoxy)phenoxy | 2,6-dichlorophenyl | 417 |
| 2-(2-propoxy)phenoxy | 2,4-dichlorophenyl | 417 |
| 2-(2-propoxy)phenoxy | 4-fluorophenyl | 366 |
| 2-(2-propoxy)phenoxy | 4-butylphenyl | 404 |
| 2-(2-propoxy)phenoxy | 2-methylphenyl | 362 |
| 2-(2-propoxy)phenoxy | phenyl | 348 |
| 2-(2-propoxy)phenoxy | 4-ethylphenyl | 376 |
| 2-(2-propoxy)phenoxy | 2,3-difluorophenyl | 384 |
| 2-(2-propoxy)phenoxy | 2,6-dimethoxyphenyl | 408 |
| 2-(2-propoxy)phenoxy | 3,4-difluorophenyl | 384 |
| 2-(2-propoxy)phenoxy | 2,5-difluorophenyl | 384 |
| 2-(2-propoxy)phenoxy | 4-ethoxyphenyl | 392 |
| 2-(2-propoxy)phenoxy | 2,4,6-trichlorophenyl | 452 |
| 2-(2-propoxy)phenoxy | 3-methylphenyl | 362 |
| 2-(2-propoxy)phenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 434 |
| 2-(2-propoxy)phenoxy | 3-methoxyphenyl | 378 |
| 2-(2-propoxy)phenoxy | 2-bromophenyl | 427 |

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| 2-(2-propoxy)phenoxy | 4-bromophenyl | 427 |
| 2-(2-propoxy)phenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 434 |
| 2-(2-propoxy)phenoxy | 3-(trifluoromethoxy)phenyl | 432 |
| 2-(2-propoxy)phenoxy | 9-fluorenon-4-yl | 450 |
| 2-(2-propoxy)phenoxy | isoxazol-5-yl | 339 |
| 2-(2-propoxy)phenoxy | benzofuroxan-5-yl | 406 |
| 2-(2-propoxy)phenoxy | 2-chloropyrid-3-yl | 384 |
| 2-(2-propoxy)phenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 455 |
| 2-(2-propoxy)phenoxy | pyridin-4-yl | 349 |
| 2-(2-propoxy)phenoxy | anthraquinon-2-yl | 478 |
| 2-(2-propoxy)phenoxy | 2-iodophenyl | 474 |
| 2-(2-propoxy)phenoxy | 4-pentylphenyl | 419 |
| 2-(2-propoxy)phenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 492 |
| 2-(2-propoxy)phenoxy | 2,6-dimethylphenyl | 376 |
| 2-(2-propoxy)phenoxy | 2,5-dimethoxyphenyl | 408 |
| 2-(2-propoxy)phenoxy | 2,5-dichloropyridin-3-yl | 418 |
| 2-(2-propoxy)phenoxy | 2-chloro-6-methoxypyridin-4-yl | 414 |
| 2-(2-propoxy)phenoxy | 2,3-dichloropyridin-5-yl | 418 |
| 2-(2-propoxy)phenoxy | 1-naphthyl | 412 |
| 2-(2-propoxy)phenoxy | 2,4-dimethoxyphenyl | 408 |
| 2-(2-propoxy)phenoxy | 3,5-bis(trifluoromethyl)phenyl | 484 |
| 2-(2-propoxy)phenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 476 |
| 2-(2-propoxy)phenoxy | pentafluorophenyl | 438 |

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| 4-fluorophenoxy | 3,4-dimethoxyphenyl | 368 |
| 4-fluorophenoxy | 2-(trifluoromethyl)phenyl | 376 |
| 4-fluorophenoxy | 2,4-difluorophenyl | 344 |
| 4-fluorophenoxy | 3-(trifluoromethyl)phenyl | 376 |
| 4-fluorophenoxy | 2-naphthyl | 358 |
| 4-fluorophenoxy | 2-methoxyphenyl | 338 |
| 4-fluorophenoxy | 3,4,5-trimethylphenyl | 398 |
| 4-fluorophenoxy | 3,4-dichlorophenyl | 377 |
| 4-fluorophenoxy | 3-bromophenyl | 387 |
| 4-fluorophenoxy | 3-pyridyl | 309 |
| 4-fluorophenoxy | 2-ethoxynaphth-1-yl | 402 |
| 4-fluorophenoxy | 2,3-dichlorophenyl | 377 |
| 4-fluorophenoxy | 6-chloropyrid-3-yl | 344 |
| 4-fluorophenoxy | 4-(trifluoromethoxy)phenyl | 392 |
| 4-fluorophenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 394 |
| 4-fluorophenoxy | 3-bromothieryl | 393 |
| 4-fluorophenoxy | 2-acetoxyphenyl | 366 |
| 4-fluorophenoxy | 5-methylisoxazol-3-yl | 313 |
| 4-fluorophenoxy | 2-(phenylthio)pyrid-3-yl | 417 |
| 4-fluorophenoxy | 2-(trifluoromethoxy)phenyl | 392 |
| 4-fluorophenoxy | 1-phenyl-5-propylpyrazin-4-yl | 416 |
| 4-fluorophenoxy | 2-ethoxyphenyl | 352 |
| 4-fluorophenoxy | 3-chlorothieryl-2-yl | 349 |
| 4-fluorophenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 368 |
| 4-fluorophenoxy | 3,5-dichlorophenyl | 377 |

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| 4-fluorophenoxy | 2-(propylthio)pyridin-3-yl | 383 |
| 4-fluorophenoxy | 2-(ethylthio)pyridin-3-yl | 369 |
| 4-fluorophenoxy | 3-bromopyridin-5-yl | 388 |
| 4-fluorophenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 330 |
| 4-fluorophenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 368 |
| 4-fluorophenoxy | 3-chlorobenzo[b]thiophen-2-yl | 399 |
| 4-fluorophenoxy | 4-chlorophenyl | 343 |
| 4-fluorophenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 389 |
| 4-fluorophenoxy | benzo[b]thiophen-2-yl | 364 |
| 4-fluorophenoxy | 3,4-dimethylphenyl | 336 |
| 4-fluorophenoxy | 2-(phenoxy)pyridin-3-yl | 401 |
| 4-fluorophenoxy | 2-(methylthio)pyridin-3-yl | 355 |
| 4-fluorophenoxy | 5-methyl-3-phenylisoxazol-4-yl | 389 |
| 4-fluorophenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 412 |
| 4-fluorophenoxy | 2-chloro-6-methylpyridin-4-yl | 358 |
| 4-fluorophenoxy | 3,5-dimethylisoxazol-4-yl | 327 |
| 4-fluorophenoxy | 1-naphthyl | 358 |
| 4-fluorophenoxy | 2-fluorophenyl | 326 |
| 4-fluorophenoxy | 4-propylphenyl | 350 |
| 4-fluorophenoxy | 3-fluorophenyl | 326 |
| 4-fluorophenoxy | 2,6-difluorophenyl | 344 |
| 4-fluorophenoxy | 2-chlorophenyl | 343 |

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| 4-fluorophenoxy | 3-(chloromethyl)phenyl | 357 |
| 4-fluorophenoxy | 4-(2-(2-methyl)propyl)phenyl | 364 |
| 4-fluorophenoxy | 3-chlorophenyl | 343 |
| 4-fluorophenoxy | 3,5-dimethoxyphenyl | 368 |
| 4-fluorophenoxy | 2,6-dichlorophenyl | 377 |
| 4-fluorophenoxy | 2,4-dichlorophenyl | 377 |
| 4-fluorophenoxy | 4-fluorophenyl | 326 |
| 4-fluorophenoxy | 4-butylphenyl | 364 |
| 4-fluorophenoxy | 2-methylphenyl | 322 |
| 4-fluorophenoxy | phenyl | 308 |
| 4-fluorophenoxy | 4-ethylphenyl | 336 |
| 4-fluorophenoxy | 2,3-difluorophenyl | 344 |
| 4-fluorophenoxy | 2,6-dimethoxyphenyl | 368 |
| 4-fluorophenoxy | 3,4-difluorophenyl | 344 |
| 4-fluorophenoxy | 2,5-difluorophenyl | 344 |
| 4-fluorophenoxy | 4-ethoxyphenyl | 352 |
| 4-fluorophenoxy | 2,4,6-trichlorophenyl | 412 |
| 4-fluorophenoxy | 3-methylphenyl | 322 |
| 4-fluorophenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 394 |
| 4-fluorophenoxy | 3-methoxyphenyl | 338 |
| 4-fluorophenoxy | 2-bromophenyl | 387 |
| 4-fluorophenoxy | 4-bromophenyl | 387 |
| 4-fluorophenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 394 |
| 4-fluorophenoxy | 3-(trifluoromethoxy)phenyl | 392 |
| 4-fluorophenoxy | 9-fluorenon-4-yl | 410 |

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| 4-fluorophenoxy | isoxazol-5-yl | 299 |
| 4-fluorophenoxy | benzofuroxan-5-yl | 366 |
| 4-fluorophenoxy | 2-chloropyrid-3-yl | 344 |
| 4-fluorophenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 415 |
| 4-fluorophenoxy | pyridin-4-yl | 309 |
| 4-fluorophenoxy | anthraquinon-2-yl | 438 |
| 4-fluorophenoxy | 2-iodophenyl | 434 |
| 4-fluorophenoxy | 4-pentylphenyl | 378 |
| 4-fluorophenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 452 |
| 4-fluorophenoxy | 2,6-dimethylphenyl | 336 |
| 4-fluorophenoxy | 2,5-dimethoxyphenyl | 368 |
| 4-fluorophenoxy | 2,5-dichloropyridin-3-yl | 378 |
| 4-fluorophenoxy | 2-chloro-6-methoxypyridin-4-yl | 374 |
| 4-fluorophenoxy | 2,3-dichloropyridin-5-yl | 378 |
| 4-fluorophenoxy | 1-naphthyl | 372 |
| 4-fluorophenoxy | 2,4-dimethoxyphenyl | 368 |
| 4-fluorophenoxy | 3,5-bis(trifluoromethyl)phenyl | 444 |
| 4-fluorophenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 436 |
| 4-fluorophenoxy | pentafluorophenyl | 398 |
| 4-chlorophenoxy | 3,4-dimethoxyphenyl | 385 |
| 4-chlorophenoxy | 2-(trifluoromethyl)phenyl | 393 |
| 4-chlorophenoxy | 2,4-difluorophenyl | 361 |
| 4-chlorophenoxy | 3-(trifluoromethyl)phenyl | 393 |
| 4-chlorophenoxy | 2-naphthyl | 375 |

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| 4-chlorophenoxy | 2-methoxyphenyl | 355 |
| 4-chlorophenoxy | 3,4,5-trimethylphenyl | 415 |
| 4-chlorophenoxy | 3,4-dichlorophenyl | 394 |
| 4-chlorophenoxy | 3-bromophenyl | 404 |
| 4-chlorophenoxy | 3-pyridyl | 326 |
| 4-chlorophenoxy | 2-ethoxynaphth-1-yl | 419 |
| 4-chlorophenoxy | 2,3-dichlorophenyl | 394 |
| 4-chlorophenoxy | 6-chloropyrid-3-yl | 360 |
| 4-chlorophenoxy | 4-(trifluoromethoxy)phenyl | 409 |
| 4-chlorophenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 411 |
| 4-chlorophenoxy | 3-bromothienyl | 410 |
| 4-chlorophenoxy | 2-acetoxyphenyl | 383 |
| 4-chlorophenoxy | 5-methylisoxazol-3-yl | 330 |
| 4-chlorophenoxy | 2-(phenylthio)pyrid-3-yl | 434 |
| 4-chlorophenoxy | 2-(trifluoromethoxy)phenyl | 409 |
| 4-chlorophenoxy | 1-phenyl-5-propylpyrazin-4-yl | 433 |
| 4-chlorophenoxy | 2-ethoxyphenyl | 369 |
| 4-chlorophenoxy | 3-chlorothien-2-yl | 365 |
| 4-chlorophenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 385 |
| 4-chlorophenoxy | 3,5-dichlorophenyl | 394 |
| 4-chlorophenoxy | 2-(propylthio)pyridin-3-yl | 400 |
| 4-chlorophenoxy | 2-(ethylthio)pyridin-3-yl | 386 |
| 4-chlorophenoxy | 3-bromopyridin-5-yl | 405 |
| 4-chlorophenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 347 |

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| 4-chlorophenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 385 |
| 4-chlorophenoxy | 3-chlorobenzo[b]thiophen-2-yl | 415 |
| 4-chlorophenoxy | 4-chlorophenyl | 359 |
| 4-chlorophenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 406 |
| 4-chlorophenoxy | benzo[b]thiophen-2-yl | 381 |
| 4-chlorophenoxy | 3,4-dimethylphenyl | 353 |
| 4-chlorophenoxy | 2-(phenoxy)pyridin-3-yl | 418 |
| 4-chlorophenoxy | 2-(methylthio)pyridin-3-yl | 372 |
| 4-chlorophenoxy | 5-methyl-3-phenylisoxazol-4-yl | 406 |
| 4-chlorophenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 428 |
| 4-chlorophenoxy | 2-chloro-6-methylpyridin-4-yl | 374 |
| 4-chlorophenoxy | 3,5-dimethylisoxazol-4-yl | 344 |
| 4-chlorophenoxy | 1-naphthyl | 375 |
| 4-chlorophenoxy | 2-fluorophenyl | 343 |
| 4-chlorophenoxy | 4-propylphenyl | 367 |
| 4-chlorophenoxy | 3-fluorophenyl | 343 |
| 4-chlorophenoxy | 2,6-difluorophenyl | 361 |
| 4-chlorophenoxy | 2-chlorophenyl | 359 |
| 4-chlorophenoxy | 3-(chloromethyl)phenyl | 373 |
| 4-chlorophenoxy | 4-(2-(2-methyl)propyl)phenyl | 381 |
| 4-chlorophenoxy | 3-chlorophenyl | 359 |

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| 4-chlorophenoxy | 3,5-dimethoxyphenyl | 385 |
| 4-chlorophenoxy | 2,6-dichlorophenyl | 394 |
| 4-chlorophenoxy | 2,4-dichlorophenyl | 394 |
| 4-chlorophenoxy | 4-fluorophenyl | 343 |
| 4-chlorophenoxy | 4-butylphenyl | 381 |
| 4-chlorophenoxy | 2-methylphenyl | 339 |
| 4-chlorophenoxy | phenyl | 325 |
| 4-chlorophenoxy | 4-ethylphenyl | 353 |
| 4-chlorophenoxy | 2,3-difluorophenyl | 361 |
| 4-chlorophenoxy | 2,6-dimethoxyphenyl | 385 |
| 4-chlorophenoxy | 3,4-difluorophenyl | 361 |
| 4-chlorophenoxy | 2,5-difluorophenyl | 361 |
| 4-chlorophenoxy | 4-ethoxyphenyl | 369 |
| 4-chlorophenoxy | 2,4,6-trichlorophenyl | 428 |
| 4-chlorophenoxy | 3-methylphenyl | 339 |
| 4-chlorophenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 411 |
| 4-chlorophenoxy | 3-methoxyphenyl | 355 |
| 4-chlorophenoxy | 2-bromophenyl | 404 |
| 4-chlorophenoxy | 4-bromophenyl | 404 |
| 4-chlorophenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 411 |
| 4-chlorophenoxy | 3-(trifluoromethoxy)phenyl | 409 |
| 4-chlorophenoxy | 9-fluorenon-4-yl | 427 |
| 4-chlorophenoxy | isoxazol-5-yl | 316 |
| 4-chlorophenoxy | benzofuroxan-5-yl | 383 |
| 4-chlorophenoxy | 2-chloropyrid-3-yl | 360 |

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| 4-chlorophenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 432 |
| 4-chlorophenoxy | pyridin-4-yl | 326 |
| 4-chlorophenoxy | anthraquinon-2-yl | 455 |
| 4-chlorophenoxy | 2-iodophenyl | 451 |
| 4-chlorophenoxy | 4-pentylphenyl | 395 |
| 4-chlorophenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 468 |
| 4-chlorophenoxy | 2,6-dimethylphenyl | 353 |
| 4-chlorophenoxy | 2,5-dimethoxyphenyl | 385 |
| 4-chlorophenoxy | 2,5-dichloropyridin-3-yl | 395 |
| 4-chlorophenoxy | 2-chloro-6-methoxypyridin-4-yl | 390 |
| 4-chlorophenoxy | 2,3-dichloropyridin-5-yl | 395 |
| 4-chlorophenoxy | 1-naphthyl | 389 |
| 4-chlorophenoxy | 2,4-dimethoxyphenyl | 385 |
| 4-chlorophenoxy | 3,5-bis(trifluoromethyl)phenyl | 461 |
| 4-chlorophenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 452 |
| 4-chlorophenoxy | pentafluorophenyl | 415 |
| 2,4-difluorophenoxy | 3,4-dimethoxyphenyl | 386 |
| 2,4-difluorophenoxy | 2-(trifluoromethyl)phenyl | 394 |
| 2,4-difluorophenoxy | 2,4-difluorophenyl | 362 |
| 2,4-difluorophenoxy | 3-(trifluoromethyl)phenyl | 394 |
| 2,4-difluorophenoxy | 2-naphthyl | 376 |
| 2,4-difluorophenoxy | 2-methoxyphenyl | 356 |
| 2,4-difluorophenoxy | 3,4,5-trimethylphenyl | 416 |

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| 2,4-difluorophenoxy | 3,4-dichlorophenyl | 395 |
| 2,4-difluorophenoxy | 3-bromophenyl | 405 |
| 2,4-difluorophenoxy | 3-pyridyl | 327 |
| 2,4-difluorophenoxy | 2-ethoxynaphth-1-yl | 420 |
| 2,4-difluorophenoxy | 2,3-dichlorophenyl | 395 |
| 2,4-difluorophenoxy | 6-chloropyrid-3-yl | 362 |
| 2,4-difluorophenoxy | 4-(trifluoromethoxy)phenyl | 410 |
| 2,4-difluorophenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 412 |
| 2,4-difluorophenoxy | 3-bromothienyl | 411 |
| 2,4-difluorophenoxy | 2-acetoxyphenyl | 384 |
| 2,4-difluorophenoxy | 5-methylisoxazol-3-yl | 331 |
| 2,4-difluorophenoxy | 2-(phenylthio)pyrid-3-yl | 435 |
| 2,4-difluorophenoxy | 2-(trifluoromethoxy)phenyl | 410 |
| 2,4-difluorophenoxy | 1-phenyl-5-propylpyrazin-4-yl | 434 |
| 2,4-difluorophenoxy | 2-ethoxyphenyl | 370 |
| 2,4-difluorophenoxy | 3-chlorothien-2-yl | 367 |
| 2,4-difluorophenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 386 |
| 2,4-difluorophenoxy | 3,5-dichlorophenyl | 395 |
| 2,4-difluorophenoxy | 2-(propylthio)pyridin-3-yl | 401 |
| 2,4-difluorophenoxy | 2-(ethylthio)pyridin-3-yl | 387 |
| 2,4-difluorophenoxy | 3-bromopyridin-5-yl | 406 |
| 2,4-difluorophenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 348 |
| 2,4-difluorophenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 386 |

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| 2,4-difluorophenoxy | 3-chlorobenzo[b]thiophen-2-yl | 417 |
| 2,4-difluorophenoxy | 4-chlorophenyl | 361 |
| 2,4-difluorophenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 407 |
| 2,4-difluorophenoxy | benzo[b]thiophen-2-yl | 382 |
| 2,4-difluorophenoxy | 3,4-dimethylphenyl | 354 |
| 2,4-difluorophenoxy | 2-(phenoxy)pyridin-3-yl | 409 |
| 2,4-difluorophenoxy | 2-(methylthio)pyridin-3-yl | 373 |
| 2,4-difluorophenoxy | 5-methyl-3-phenylisoxazol-4-yl | 407 |
| 2,4-difluorophenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 430 |
| 2,4-difluorophenoxy | 2-chloro-6-methylpyridin-4-yl | 376 |
| 2,4-difluorophenoxy | 3,5-dimethylisoxazol-4-yl | 345 |
| 2,4-difluorophenoxy | 1-naphthyl | 376 |
| 2,4-difluorophenoxy | 2-fluorophenyl | 344 |
| 2,4-difluorophenoxy | 4-propylphenyl | 368 |
| 2,4-difluorophenoxy | 3-fluorophenyl | 344 |
| 2,4-difluorophenoxy | 2,6-difluorophenyl | 362 |
| 2,4-difluorophenoxy | 2-chlorophenyl | 361 |
| 2,4-difluorophenoxy | 3-(chloromethyl)phenyl | 375 |
| 2,4-difluorophenoxy | 4-(2-(2-methyl)propyl)phenyl | 382 |
| 2,4-difluorophenoxy | 3-chlorophenyl | 361 |
| 2,4-difluorophenoxy | 3,5-dimethoxyphenyl | 386 |
| 2,4-difluorophenoxy | 2,6-dichlorophenyl | 395 |

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| 2,4-difluorophenoxy | 2,4-dichlorophenyl | 392 |
| 2,4-difluorophenoxy | 4-fluorophenyl | 344 |
| 2,4-difluorophenoxy | 4-butylphenyl | 382 |
| 2,4-difluorophenoxy | 2-methylphenyl | 340 |
| 2,4-difluorophenoxy | phenyl | 326 |
| 2,4-difluorophenoxy | 4-ethylphenyl | 354 |
| 2,4-difluorophenoxy | 2,3-difluorophenyl | 362 |
| 2,4-difluorophenoxy | 2,6-dimethoxyphenyl | 386 |
| 2,4-difluorophenoxy | 3,4-difluorophenyl | 362 |
| 2,4-difluorophenoxy | 2,5-difluorophenyl | 362 |
| 2,4-difluorophenoxy | 4-ethoxyphenyl | 370 |
| 2,4-difluorophenoxy | 2,4,6-trichlorophenyl | 430 |
| 2,4-difluorophenoxy | 3-methylphenyl | 340 |
| 2,4-difluorophenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 412 |
| 2,4-difluorophenoxy | 3-methoxyphenyl | 356 |
| 2,4-difluorophenoxy | 2-bromophenyl | 405 |
| 2,4-difluorophenoxy | 4-bromophenyl | 405 |
| 2,4-difluorophenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 412 |
| 2,4-difluorophenoxy | 3-(trifluoromethoxy)phenyl | 410 |
| 2,4-difluorophenoxy | 9-fluorenon-4-yl | 428 |
| 2,4-difluorophenoxy | isoxazol-5-yl | 317 |
| 2,4-difluorophenoxy | benzofuroxan-5-yl | 384 |
| 2,4-difluorophenoxy | 2-chloropyrid-3-yl | 362 |
| 2,4-difluorophenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 433 |
| 2,4-difluorophenoxy | pyridin-4-yl | 327 |

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| 2,4-difluorophenoxy | anthraquinon-2-yl | 456 |
| 2,4-difluorophenoxy | 2-iodophenyl | 452 |
| 2,4-difluorophenoxy | 4-pentylphenyl | 396 |
| 2,4-difluorophenoxy | 2-(4-chlorophenylthio) pyridin-3-yl | 470 |
| 2,4-difluorophenoxy | 2,6-dimethylphenyl | 354 |
| 2,4-difluorophenoxy | 2,5-dimethoxyphenyl | 386 |
| 2,4-difluorophenoxy | 2,5-dichloropyridin-3-yl | 396 |
| 2,4-difluorophenoxy | 2-chloro-6-methoxypyridin- 4-yl | 392 |
| 2,4-difluorophenoxy | 2,3-dichloropyridin-5-yl | 396 |
| 2,4-difluorophenoxy | 1-naphthyl | 390 |
| 2,4-difluorophenoxy | 2,4-dimethoxyphenyl | 386 |
| 2,4-difluorophenoxy | 3,5- bis(trifluoromethyl)phenyl | 462 |
| 2,4-difluorophenoxy | 2-(4- chlorophenoxy)pyridin-3-yl | 454 |
| 2,4-difluorophenoxy | pentafluorophenyl | 416 |
| 4-thiomethylphenoxy | 3,4-dimethoxyphenyl | 396 |
| 4-thiomethylphenoxy | 2-(trifluoromethyl)phenyl | 404 |
| 4-thiomethylphenoxy | 2,4-difluorophenyl | 372 |
| 4-thiomethylphenoxy | 3-(trifluoromethyl)phenyl | 404 |
| 4-thiomethylphenoxy | 2-naphthyl | 386 |
| 4-thiomethylphenoxy | 2-methoxyphenyl | 366 |
| 4-thiomethylphenoxy | 3,4,5,-trimethylphenyl | 426 |
| 4-thiomethylphenoxy | 3,4-dichlorophenyl | 405 |
| 4-thiomethylphenoxy | 3-bromophenyl | 415 |
| 4-thiomethylphenoxy | 3-pyridyl | 337 |
| 4-thiomethylphenoxy | 2-ethoxynaphth-1-yl | 430 |

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| 4-thiomethylphenoxy | 2,3-dichlorophenyl | 405 |
| 4-thiomethylphenoxy | 6-chloropyrid-3-yl | 372 |
| 4-thiomethylphenoxy | 4-(trifluoromethoxy)phenyl | 420 |
| 4-thiomethylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 422 |
| 4-thiomethylphenoxy | 3-bromothieryl | 421 |
| 4-thiomethylphenoxy | 2-acetoxyphenyl | 394 |
| 4-thiomethylphenoxy | 5-methylisoxazol-3-yl | 341 |
| 4-thiomethylphenoxy | 2-(phenylthio)pyrid-3-yl | 446 |
| 4-thiomethylphenoxy | 2-(trifluoromethoxy)phenyl | 420 |
| 4-thiomethylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 445 |
| 4-thiomethylphenoxy | 2-ethoxyphenyl | 380 |
| 4-thiomethylphenoxy | 3-chlorothien-2-yl | 377 |
| 4-thiomethylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 396 |
| 4-thiomethylphenoxy | 3,5-dichlorophenyl | 405 |
| 4-thiomethylphenoxy | 2-(propylthio)pyridin-3-yl | 412 |
| 4-thiomethylphenoxy | 2-(ethylthio)pyridin-3-yl | 397 |
| 4-thiomethylphenoxy | 3-bromopyridin-5-yl | 416 |
| 4-thiomethylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 358 |
| 4-thiomethylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 396 |
| 4-thiomethylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 427 |
| 4-thiomethylphenoxy | 4-chlorophenyl | 371 |
| 4-thiomethylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 417 |

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| 4-thiomethylphenoxy | benzo[b]thiophen-2-yl | 392 |
| 4-thiomethylphenoxy | 3,4-dimethylphenyl | 364 |
| 4-thiomethylphenoxy | 2-(phenoxy)pyridin-3-yl | 429 |
| 4-thiomethylphenoxy | 2-(methylthio)pyridin-3-yl | 383 |
| 4-thiomethylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 417 |
| 4-thiomethylphenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 440 |
| 4-thiomethylphenoxy | 2-chloro-6-methylpyridin-4-yl | 386 |
| 4-thiomethylphenoxy | 3,5-dimethylisoxazol-4-yl | 355 |
| 4-thiomethylphenoxy | 1-naphthyl | 386 |
| 4-thiomethylphenoxy | 2-fluorophenyl | 354 |
| 4-thiomethylphenoxy | 4-propylphenyl | 378 |
| 4-thiomethylphenoxy | 3-fluorophenyl | 354 |
| 4-thiomethylphenoxy | 2,6-difluorophenyl | 372 |
| 4-thiomethylphenoxy | 2-chlorophenyl | 371 |
| 4-thiomethylphenoxy | 3-(chloromethyl)phenyl | 385 |
| 4-thiomethylphenoxy | 4-(2-(2-methyl)propyl)phenyl | 392 |
| 4-thiomethylphenoxy | 3-chlorophenyl | 371 |
| 4-thiomethylphenoxy | 3,5-dimethoxyphenyl | 396 |
| 4-thiomethylphenoxy | 2,6-dichlorophenyl | 405 |
| 4-thiomethylphenoxy | 2,4-dichlorophenyl | 405 |
| 4-thiomethylphenoxy | 4-fluorophenyl | 354 |
| 4-thiomethylphenoxy | 4-butylphenyl | 392 |
| 4-thiomethylphenoxy | 2-methylphenyl | 350 |
| 4-thiomethylphenoxy | phenyl | 336 |

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| 4-thiomethylphenoxy | 4-ethylphenyl | 364 |
| 4-thiomethylphenoxy | 2,3-difluorophenyl | 372 |
| 4-thiomethylphenoxy | 2,6-dimethoxyphenyl | 396 |
| 4-thiomethylphenoxy | 3,4-difluorophenyl | 372 |
| 4-thiomethylphenoxy | 2,5-difluorophenyl | 372 |
| 4-thiomethylphenoxy | 4-ethoxyphenyl | 380 |
| 4-thiomethylphenoxy | 2,4,6-trichlorophenyl | 440 |
| 4-thiomethylphenoxy | 3-methylphenyl | 350 |
| 4-thiomethylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 422 |
| 4-thiomethylphenoxy | 3-methoxyphenyl | 366 |
| 4-thiomethylphenoxy | 2-bromophenyl | 415 |
| 4-thiomethylphenoxy | 4-bromophenyl | 415 |
| 4-thiomethylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 422 |
| 4-thiomethylphenoxy | 3-(trifluoromethoxy)phenyl | 420 |
| 4-thiomethylphenoxy | 9-fluorenon-4-yl | 438 |
| 4-thiomethylphenoxy | isoxazol-5-yl | 327 |
| 4-thiomethylphenoxy | benzofuroxan-5-yl | 394 |
| 4-thiomethylphenoxy | 2-chloropyrid-3-yl | 372 |
| 4-thiomethylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 443 |
| 4-thiomethylphenoxy | pyridin-4-yl | 337 |
| 4-thiomethylphenoxy | anthraquinon-2-yl | 466 |
| 4-thiomethylphenoxy | 2-iodophenyl | 462 |
| 4-thiomethylphenoxy | 4-pentylphenyl | 407 |
| 4-thiomethylphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 480 |
| 4-thiomethylphenoxy | 2,6-dimethylphenyl | 364 |

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| 4-thiomethylphenoxy | 2,5-dimethoxyphenyl | 396 |
| 4-thiomethylphenoxy | 2,5-dichloropyridin-3-yl | 406 |
| 4-thiomethylphenoxy | 2-chloro-6-methoxypyridin-4-yl | 402 |
| 4-thiomethylphenoxy | 2,3-dichloropyridin-5-yl | 406 |
| 4-thiomethylphenoxy | 1-naphthyl | 400 |
| 4-thiomethylphenoxy | 2,4-dimethoxyphenyl | 396 |
| 4-thiomethylphenoxy | 3,5-bis(trifluoromethyl)phenyl | 372 |
| 4-thiomethylphenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 464 |
| 4-thiomethylphenoxy | pentafluorophenyl | 426 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,4-dimethoxyphenyl | 406 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(trifluoromethyl)phenyl | 414 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,4-difluorophenyl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-(trifluoromethyl)phenyl | 414 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-naphthyl | 396 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-methoxyphenyl | 376 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,4,5-trimethylphenyl | 436 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,4-dichlorophenyl | 415 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-bromophenyl | 425 |

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| 4-(2-(2-methyl)propyl)phenoxy | 3-pyridyl | 347 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-ethoxynaphth-1-yl | 441 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,3-dichlorophenyl | 415 |
| 4-(2-(2-methyl)propyl)phenoxy | 6-chloropyrid-3-yl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-(trifluoromethoxy)phenyl | 430 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 432 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-bromothieryl | 431 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-acetoxyphenyl | 404 |
| 4-(2-(2-methyl)propyl)phenoxy | 5-methylisoxazol-3-yl | 351 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(phenylthio)pyrid-3-yl | 456 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(trifluoromethoxy)phenyl | 430 |
| 4-(2-(2-methyl)propyl)phenoxy | 1-phenyl-5-propylpyrazin-4-yl | 455 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-ethoxyphenyl | 390 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-chlorothieryl-2-yl | 387 |
| 4-(2-(2-methyl)propyl)phenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 406 |

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| 4-(2-(2-methyl)propyl)phenoxy | 3,5-dichlorophenyl | 415 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(propylthio)pyridin-3-yl | 422 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(ethylthio)pyridin-3-yl | 407 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-bromopyridin-5-yl | 426 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 368 |
| 4-(2-(2-methyl)propyl)phenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 406 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-chlorobenzo[b]thiophen-2-yl | 437 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-chlorophenyl | 381 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 427 |
| 4-(2-(2-methyl)propyl)phenoxy | benzo[b]thiophen-2-yl | 402 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,4-dimethylphenyl | 374 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(phenoxy)pyridin-3-yl | 439 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(methylthio)pyridin-3-yl | 393 |
| 4-(2-(2-methyl)propyl)phenoxy | 5-methyl-3-phenylisoxazol-4-yl | 427 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 450 |

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| 4-(2-(2-methyl)propyl)phenoxy | 2-chloro-6-methylpyridin-4-yl | 396 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,5-dimethylisoxazol-4-yl | 365 |
| 4-(2-(2-methyl)propyl)phenoxy | 1-naphthyl | 396 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-fluorophenyl | 364 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-propylphenyl | 388 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-fluorophenyl | 364 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,6-difluorophenyl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-chlorophenyl | 381 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-(chloromethyl)phenyl | 395 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-(2-(2-methyl)propyl)phenyl | 402 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-chlorophenyl | 381 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,5-dimethoxyphenyl | 406 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,6-dichlorophenyl | 415 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,4-dichlorophenyl | 415 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-fluorophenyl | 364 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-butylphenyl | 402 |

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| 4-(2-(2-methyl)propyl)phenoxy | 2-methylphenyl | 360 |
| 4-(2-(2-methyl)propyl)phenoxy | phenyl | 346 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-ethylphenyl | 374 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,3-difluorophenyl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,6-dimethoxyphenyl | 406 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,4-difluorophenyl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,5-difluorophenyl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-ethoxyphenyl | 390 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,4,6-trichlorophenyl | 450 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-methylphenyl | 360 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 432 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-methoxyphenyl | 376 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-bromophenyl | 425 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-bromophenyl | 425 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 432 |
| 4-(2-(2-methyl)propyl)phenoxy | 3-(trifluoromethoxy)phenyl | 430 |

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| 4-(2-(2-methyl)propyl)phenoxy | 9-fluorenon-4-yl | 448 |
| 4-(2-(2-methyl)propyl)phenoxy | isoxazol-5-yl | 338 |
| 4-(2-(2-methyl)propyl)phenoxy | benzofuroxan-5-yl | 404 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-chloropyrid-3-yl | 382 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 454 |
| 4-(2-(2-methyl)propyl)phenoxy | pyridin-4-yl | 347 |
| 4-(2-(2-methyl)propyl)phenoxy | anthraquinon-2-yl | 476 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-iodophenyl | 472 |
| 4-(2-(2-methyl)propyl)phenoxy | 4-pentylphenyl | 417 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 490 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,6-dimethylphenyl | 374 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,5-dimethoxyphenyl | 406 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,5-dichloropyridin-3-yl | 416 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-chloro-6-methoxypyridin-4-yl | 412 |
| 4-(2-(2-methyl)propyl)phenoxy | 2,3-dichloropyridin-5-yl | 416 |
| 4-(2-(2-methyl)propyl)phenoxy | 1-naphthyl | 410 |

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| 4-(2-(2-methyl)propyl)phenoxy | 2,4-dimethoxyphenyl | 406 |
| 4-(2-(2-methyl)propyl)phenoxy | 3,5-bis(trifluoromethyl)phenyl | 482 |
| 4-(2-(2-methyl)propyl)phenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 474 |
| 4-(2-(2-methyl)propyl)phenoxy | pentafluorophenyl | 436 |
| 2,3-dimethylphenoxy | 3,4-dimethoxyphenyl | 378 |
| 2,3-dimethylphenoxy | 2-(trifluoromethyl)phenyl | 386 |
| 2,3-dimethylphenoxy | 2,4-difluorophenyl | 354 |
| 2,3-dimethylphenoxy | 3-(trifluoromethyl)phenyl | 386 |
| 2,3-dimethylphenoxy | 2-naphthyl | 368 |
| 2,3-dimethylphenoxy | 2-methoxyphenyl | 348 |
| 2,3-dimethylphenoxy | 3,4,5-trimethylphenyl | 408 |
| 2,3-dimethylphenoxy | 3,4-dichlorophenyl | 387 |
| 2,3-dimethylphenoxy | 3-bromophenyl | 397 |
| 2,3-dimethylphenoxy | 3-pyridyl | 319 |
| 2,3-dimethylphenoxy | 2-ethoxynaphth-1-yl | 412 |
| 2,3-dimethylphenoxy | 2,3-dichlorophenyl | 387 |
| 2,3-dimethylphenoxy | 6-chloropyrid-3-yl | 354 |
| 2,3-dimethylphenoxy | 4-(trifluoromethoxy)phenyl | 402 |
| 2,3-dimethylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 404 |
| 2,3-dimethylphenoxy | 3-bromothieryl | 403 |
| 2,3-dimethylphenoxy | 2-acetoxyphenyl | 376 |
| 2,3-dimethylphenoxy | 5-methylisoxazol-3-yl | 323 |
| 2,3-dimethylphenoxy | 2-(phenylthio)pyrid-3-yl | 427 |
| 2,3-dimethylphenoxy | 2-(trifluoromethoxy)phenyl | 402 |

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| 2,3-dimethylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 426 |
| 2,3-dimethylphenoxy | 2-ethoxyphenyl | 362 |
| 2,3-dimethylphenoxy | 3-chlorothien-2-yl | 359 |
| 2,3-dimethylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 378 |
| 2,3-dimethylphenoxy | 3,5-dichlorophenyl | 387 |
| 2,3-dimethylphenoxy | 2-(propylthio)pyridin-3-yl | 393 |
| 2,3-dimethylphenoxy | 2-(ethylthio)pyridin-3-yl | 379 |
| 2,3-dimethylphenoxy | 3-bromopyridin-5-yl | 398 |
| 2,3-dimethylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 340 |
| 2,3-dimethylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 378 |
| 2,3-dimethylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 409 |
| 2,3-dimethylphenoxy | 4-chlorophenyl | 353 |
| 2,3-dimethylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 399 |
| 2,3-dimethylphenoxy | benzo[b]thiophen-2-yl | 374 |
| 2,3-dimethylphenoxy | 3,4-dimethylphenyl | 346 |
| 2,3-dimethylphenoxy | 2-(phenoxy)pyridin-3-yl | 411 |
| 2,3-dimethylphenoxy | 2-(methylthio)pyridin-3-yl | 365 |
| 2,3-dimethylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 399 |
| 2,3-dimethylphenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 422 |
| 2,3-dimethylphenoxy | 2-chloro-6-methylpyridin-4-yl | 368 |

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| 2,3-dimethylphenoxy | 3,5-dimethylisoxazol-4-yl | 337 |
| 2,3-dimethylphenoxy | 1-naphthyl | 368 |
| 2,3-dimethylphenoxy | 2-fluorophenyl | 336 |
| 2,3-dimethylphenoxy | 4-propylphenyl | 360 |
| 2,3-dimethylphenoxy | 3-fluorophenyl | 336 |
| 2,3-dimethylphenoxy | 2,6-difluorophenyl | 354 |
| 2,3-dimethylphenoxy | 2-chlorophenyl | 353 |
| 2,3-dimethylphenoxy | 3-(chloromethyl)phenyl | 368 |
| 2,3-dimethylphenoxy | 4-(2-(2-methyl)propyl) phenyl | 374 |
| 2,3-dimethylphenoxy | 3-chlorophenyl | 353 |
| 2,3-dimethylphenoxy | 3,5-dimethoxyphenyl | 378 |
| 2,3-dimethylphenoxy | 2,6-dichlorophenyl | 387 |
| 2,3-dimethylphenoxy | 2,4-dichlorophenyl | 387 |
| 2,3-dimethylphenoxy | 4-fluorophenyl | 336 |
| 2,3-dimethylphenoxy | 4-butylphenyl | 374 |
| 2,3-dimethylphenoxy | 2-methylphenyl | 332 |
| 2,3-dimethylphenoxy | phenyl | 318 |
| 2,3-dimethylphenoxy | 4-ethylphenyl | 346 |
| 2,3-dimethylphenoxy | 2,3-difluorophenyl | 354 |
| 2,3-dimethylphenoxy | 2,6-dimethoxyphenyl | 378 |
| 2,3-dimethylphenoxy | 3,4-difluorophenyl | 354 |
| 2,3-dimethylphenoxy | 2,5-difluorophenyl | 354 |
| 2,3-dimethylphenoxy | 4-ethoxyphenyl | 362 |
| 2,3-dimethylphenoxy | 2,4,6-trichlorophenyl | 422 |
| 2,3-dimethylphenoxy | 3-methylphenyl | 332 |
| 2,3-dimethylphenoxy | 2-fluoro-5- (trifluoromethyl)phenyl | 404 |

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| 2,3-dimethylphenoxy | 3-methoxyphenyl | 348 |
| 2,3-dimethylphenoxy | 2-bromophenyl | 397 |
| 2,3-dimethylphenoxy | 4-bromophenyl | 397 |
| 2,3-dimethylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 404 |
| 2,3-dimethylphenoxy | 3-(trifluoromethoxy)phenyl | 402 |
| 2,3-dimethylphenoxy | 9-fluorenon-4-yl | 420 |
| 2,3-dimethylphenoxy | isoxazol-5-yl | 609 |
| 2,3-dimethylphenoxy | benzofuroxan-5-yl | 376 |
| 2,3-dimethylphenoxy | 2-chloropyrid-3-yl | 354 |
| 2,3-dimethylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 425 |
| 2,3-dimethylphenoxy | pyridin-4-yl | 319 |
| 2,3-dimethylphenoxy | anthraquinon-2-yl | 448 |
| 2,3-dimethylphenoxy | 2-iodophenyl | 444 |
| 2,3-dimethylphenoxy | 4-pentylphenyl | 388 |
| 2,3-dimethylphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 462 |
| 2,3-dimethylphenoxy | 2,6-dimethylphenyl | 346 |
| 2,3-dimethylphenoxy | 2,5-dimethoxyphenyl | 378 |
| 2,3-dimethylphenoxy | 2,5-dichloropyridin-3-yl | 388 |
| 2,3-dimethylphenoxy | 2-chloro-6-methoxypyridin-4-yl | 384 |
| 2,3-dimethylphenoxy | 2,3-dichloropyridin-5-yl | 388 |
| 2,3-dimethylphenoxy | 1-naphthyl | 382 |
| 2,3-dimethylphenoxy | 2,4-dimethoxyphenyl | 378 |
| 2,3-dimethylphenoxy | 3,5-bis(trifluoromethyl)phenyl | 454 |

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| 2,3-dimethylphenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 446 |
| 2,3-dimethylphenoxy | pentafluorophenyl | 408 |
| 3,5-(bis-2-propyl)phenoxy | 3,4-dimethoxyphenyl | 434 |
| 3,5-(bis-2-propyl)phenoxy | 2-(trifluoromethyl)phenyl | 442 |
| 3,5-(bis-2-propyl)phenoxy | 2,4-difluorophenyl | 410 |
| 3,5-(bis-2-propyl)phenoxy | 3-(trifluoromethyl)phenyl | 442 |
| 3,5-(bis-2-propyl)phenoxy | 2-naphthyl | 425 |
| 3,5-(bis-2-propyl)phenoxy | 2-methoxyphenyl | 404 |
| 3,5-(bis-2-propyl)phenoxy | 3,4,5-trimethylphenyl | 465 |
| 3,5-(bis-2-propyl)phenoxy | 3,4-dichlorophenyl | 443 |
| 3,5-(bis-2-propyl)phenoxy | 3-bromophenyl | 453 |
| 3,5-(bis-2-propyl)phenoxy | 3-pyridyl | 375 |
| 3,5-(bis-2-propyl)phenoxy | 2-ethoxynaphth-1-yl | 469 |
| 3,5-(bis-2-propyl)phenoxy | 2,3-dichlorophenyl | 443 |
| 3,5-(bis-2-propyl)phenoxy | 6-chloropyrid-3-yl | 410 |
| 3,5-(bis-2-propyl)phenoxy | 4-(trifluoromethoxy)phenyl | 458 |

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| 3,5-(bis-2-propyl)phenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 460 |
| 3,5-(bis-2-propyl)phenoxy | 3-bromothieryl | 459 |
| 3,5-(bis-2-propyl)phenoxy | 2-acetoxyphenyl | 432 |
| 3,5-(bis-2-propyl)phenoxy | 5-methylisoxazol-3-yl | 379 |
| 3,5-(bis-2-propyl)phenoxy | 2-(phenylthio)pyrid-3-yl | 484 |
| 3,5-(bis-2-propyl)phenoxy | 2-(trifluoromethoxy)phenyl | 458 |
| 3,5-(bis-2-propyl)phenoxy | 1-phenyl-5-propylpyrazin-4-yl | 483 |
| 3,5-(bis-2-propyl)phenoxy | 2-ethoxyphenyl | 418 |
| 3,5-(bis-2-propyl)phenoxy | 3-chlorothiophen-2-yl | 415 |
| 3,5-(bis-2-propyl)phenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 435 |
| 3,5-(bis-2-propyl)phenoxy | 3,5-dichlorophenyl | 443 |
| 3,5-(bis-2-propyl)phenoxy | 2-(propylthio)pyridin-3-yl | 450 |
| 3,5-(bis-2-propyl)phenoxy | 2-(ethylthio)pyridin-3-yl | 436 |
| 3,5-(bis-2-propyl)phenoxy | 3-bromopyridin-5-yl | 454 |
| 3,5-(bis-2-propyl)phenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 396 |
| 3,5-(bis-2-propyl)phenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 435 |

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| 3,5-(bis-2-propyl)phenoxy | 3-chlorobenzo[b]thiophen-2-yl | 465 |
| 3,5-(bis-2-propyl)phenoxy | 4-chlorophenyl | 409 |
| 3,5-(bis-2-propyl)phenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 456 |
| 3,5-(bis-2-propyl)phenoxy | benzo[b]thiophen-2-yl | 431 |
| 3,5-(bis-2-propyl)phenoxy | 3,4-dimethylphenyl | 402 |
| 3,5-(bis-2-propyl)phenoxy | 2-(phenoxy)pyridin-3-yl | 468 |
| 3,5-(bis-2-propyl)phenoxy | 2-(methylthio)pyridin-3-yl | 422 |
| 3,5-(bis-2-propyl)phenoxy | 5-methyl-3-phenylisoxazol-4-yl | 456 |
| 3,5-(bis-2-propyl)phenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 478 |
| 3,5-(bis-2-propyl)phenoxy | 2-chloro-6-methylpyridin-4-yl | 424 |
| 3,5-(bis-2-propyl)phenoxy | 3,5-dimethylisoxazol-4-yl | 393 |
| 3,5-(bis-2-propyl)phenoxy | 1-naphthyl | 425 |
| 3,5-(bis-2-propyl)phenoxy | 2-fluorophenyl | 392 |
| 3,5-(bis-2-propyl)phenoxy | 4-propylphenyl | 417 |
| 3,5-(bis-2-propyl)phenoxy | 3-fluorophenyl | 392 |

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| 3,5-(bis-2-propyl)phenoxy | 2,6-difluorophenyl | 410 |
| 3,5-(bis-2-propyl)phenoxy | 2-chlorophenyl | 409 |
| 3,5-(bis-2-propyl)phenoxy | 3-(chloromethyl)phenyl | 423 |
| 3,5-(bis-2-propyl)phenoxy | 4-(2-(2-methyl)propyl)phenyl | 431 |
| 3,5-(bis-2-propyl)phenoxy | 3-chlorophenyl | 409 |
| 3,5-(bis-2-propyl)phenoxy | 3,5-dimethoxyphenyl | 434 |
| 3,5-(bis-2-propyl)phenoxy | 2,6-dichlorophenyl | 443 |
| 3,5-(bis-2-propyl)phenoxy | 2,4-dichlorophenyl | 443 |
| 3,5-(bis-2-propyl)phenoxy | 4-fluorophenyl | 392 |
| 3,5-(bis-2-propyl)phenoxy | 4-butylphenyl | 431 |
| 3,5-(bis-2-propyl)phenoxy | 2-methylphenyl | 388 |
| 3,5-(bis-2-propyl)phenoxy | phenyl | 374 |
| 3,5-(bis-2-propyl)phenoxy | 4-ethylphenyl | 402 |
| 3,5-(bis-2-propyl)phenoxy | 2,3-difluorophenyl | 410 |
| 3,5-(bis-2-propyl)phenoxy | 2,6-dimethoxyphenyl | 434 |

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| 3,5-(bis-2-propyl)phenoxy | 3,4-difluorophenyl | 410 |
| 3,5-(bis-2-propyl)phenoxy | 2,5-difluorophenyl | 410 |
| 3,5-(bis-2-propyl)phenoxy | 4-ethoxyphenyl | 418 |
| 3,5-(bis-2-propyl)phenoxy | 2,4,6-trichlorophenyl | 478 |
| 3,5-(bis-2-propyl)phenoxy | 3-methylphenyl | 388 |
| 3,5-(bis-2-propyl)phenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 460 |
| 3,5-(bis-2-propyl)phenoxy | 3-methoxyphenyl | 404 |
| 3,5-(bis-2-propyl)phenoxy | 2-bromophenyl | 453 |
| 3,5-(bis-2-propyl)phenoxy | 4-bromophenyl | 453 |
| 3,5-(bis-2-propyl)phenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 460 |
| 3,5-(bis-2-propyl)phenoxy | 3-(trifluoromethoxy)phenyl | 458 |
| 3,5-(bis-2-propyl)phenoxy | 9-fluorenon-4-yl | 477 |
| 3,5-(bis-2-propyl)phenoxy | isoxazol-5-yl | 365 |
| 3,5-(bis-2-propyl)phenoxy | benzofuroxan-5-yl | 432 |
| 3,5-(bis-2-propyl)phenoxy | 2-chloropyrid-3-yl | 410 |

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| 3,5-(bis-2-propyl)phenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 482 |
| 3,5-(bis-2-propyl)phenoxy | pyridin-4-yl | 375 |
| 3,5-(bis-2-propyl)phenoxy | anthraquinon-2-yl | 505 |
| 3,5-(bis-2-propyl)phenoxy | 2-iodophenyl | 500 |
| 3,5-(bis-2-propyl)phenoxy | 4-pentylphenyl | 445 |
| 3,5-(bis-2-propyl)phenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 518 |
| 3,5-(bis-2-propyl)phenoxy | 2,6-dimethylphenyl | 402 |
| 3,5-(bis-2-propyl)phenoxy | 2,5-dimethoxyphenyl | 434 |
| 3,5-(bis-2-propyl)phenoxy | 2,5-dichloropyridin-3-yl | 444 |
| 3,5-(bis-2-propyl)phenoxy | 2-chloro-6-methoxypyridin-4-yl | 440 |
| 3,5-(bis-2-propyl)phenoxy | 2,3-dichloropyridin-5-yl | 444 |
| 3,5-(bis-2-propyl)phenoxy | 1-naphthyl | 439 |
| 3,5-(bis-2-propyl)phenoxy | 2,4-dimethoxyphenyl | 434 |
| 3,5-(bis-2-propyl)phenoxy | 3,5-bis(trifluoromethyl)phenyl | 510 |
| 3,5-(bis-2-propyl)phenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 502 |

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| 3,5-(bis-2-propyl)phenoxy | pentafluorophenyl | 464 |
| 3-trifluoromethyl phenoxy | 3,4-dimethoxyphenyl | 418 |
| 3-trifluoromethyl phenoxy | 2-(trifluoromethyl)phenyl | 426 |
| 3-trifluoromethyl phenoxy | 2,4-difluorophenyl | 394 |
| 3-trifluoromethyl phenoxy | 3-(trifluoromethyl)phenyl | 426 |
| 3-trifluoromethyl phenoxy | 2-naphthyl | 408 |
| 3-trifluoromethyl phenoxy | 2-methoxyphenyl | 388 |
| 3-trifluoromethyl phenoxy | 3,4,5-trimethylphenyl | 448 |
| 3-trifluoromethyl phenoxy | 3,4-dichlorophenyl | 427 |
| 3-trifluoromethyl phenoxy | 3-bromophenyl | 437 |
| 3-trifluoromethyl phenoxy | 3-pyridyl | 359 |
| 3-trifluoromethyl phenoxy | 2-ethoxynaphth-1-yl | 452 |
| 3-trifluoromethyl phenoxy | 2,3-dichlorophenyl | 427 |
| 3-trifluoromethyl phenoxy | 6-chloropyrid-3-yl | 394 |
| 3-trifluoromethyl phenoxy | 4-(trifluoromethoxy)phenyl | 442 |

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| 3-trifluoromethyl phenoxy | 2-fluoro-4- (trifluoromethyl)phenyl | 444 |
| 3-trifluoromethyl phenoxy | 3-bromothienyl | 443 |
| 3-trifluoromethyl phenoxy | 2-acetoxyphenyl | 416 |
| 3-trifluoromethyl phenoxy | 5-methylisoxazol-3-yl | 363 |
| 3-trifluoromethyl phenoxy | 2-(phenylthio)pyrid-3-yl | 467 |
| 3-trifluoromethyl phenoxy | 2-(trifluoromethoxy)phenyl | 442 |
| 3-trifluoromethyl phenoxy | 1-phenyl-5-propylpyrazin- 4-yl | 466 |
| 3-trifluoromethyl phenoxy | 2-ethoxyphenyl | 402 |
| 3-trifluoromethyl phenoxy | 3-chlorothien-2-yl | 399 |
| 3-trifluoromethyl phenoxy | 1-(2-(2-methyl)propyl)-3- methylpyrazol-5-yl | 418 |
| 3-trifluoromethyl phenoxy | 3,5-dichlorophenyl | 427 |
| 3-trifluoromethyl phenoxy | 2-(propylthio)pyridin-3-yl | 433 |
| 3-trifluoromethyl phenoxy | 2-(ethylthio)pyridin-3-yl | 419 |
| 3-trifluoromethyl phenoxy | 3-bromopyridin-5-yl | 438 |
| 3-trifluoromethyl phenoxy | 4-methyl-1,2,3-thiadiazol- 5-yl | 380 |

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| 3-trifluoromethyl phenoxy | 1-methyl-3-(2-(2- methyl)propyl)pyrazol-5-yl | 418 |
| 3-trifluoromethyl phenoxy | 3-chlorobenzo[b]thiophen- 2-yl | 449 |
| 3-trifluoromethyl phenoxy | 4-chlorophenyl | 393 |
| 3-trifluoromethyl phenoxy | 4-methyl-2-phenyl-1,2,3- triazol-5-yl | 439 |
| 3-trifluoromethyl phenoxy | benzo[b]thiophen-2-yl | 414 |
| 3-trifluoromethyl phenoxy | 3,4-dimethylphenyl | 386 |
| 3-trifluoromethyl phenoxy | 2-(phenoxy)pyridin-3-yl | 451 |
| 3-trifluoromethyl phenoxy | 2-(methylthio)pyridin-3-yl | 405 |
| 3-trifluoromethyl phenoxy | 5-methyl-3-phenylisoxazol- 4-yl | 439 |
| 3-trifluoromethyl phenoxy | 4-chloro-1,3-dimethyl pyrazolo[3,4-b]pyridin-3- yl | 462 |
| 3-trifluoromethyl phenoxy | 2-chloro-6-methylpyridin- 4-yl | 408 |
| 3-trifluoromethyl phenoxy | 3,5-dimethylisoxazol-4-yl | 377 |
| 3-trifluoromethyl phenoxy | 1-naphthyl | 408 |
| 3-trifluoromethyl phenoxy | 2-fluorophenyl | 476 |
| 3-trifluoromethyl phenoxy | 4-propylphenyl | 400 |

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| 3-trifluoromethyl phenoxy | 3-fluorophenyl | 376 |
| 3-trifluoromethyl phenoxy | 2,6-difluorophenyl | 394 |
| 3-trifluoromethyl phenoxy | 2-chlorophenyl | 393 |
| 3-trifluoromethyl phenoxy | 3-(chloromethyl)phenyl | 407 |
| 3-trifluoromethyl phenoxy | 4-(2-(2-methyl)propyl) phenyl | 414 |
| 3-trifluoromethyl phenoxy | 3-chlorophenyl | 393 |
| 3-trifluoromethyl phenoxy | 3,5-dimethoxyphenyl | 418 |
| 3-trifluoromethyl phenoxy | 2,6-dichlorophenyl | 427 |
| 3-trifluoromethyl phenoxy | 2,4-dichlorophenyl | 427 |
| 3-trifluoromethyl phenoxy | 4-fluorophenyl | 376 |
| 3-trifluoromethyl phenoxy | 4-butylphenyl | 414 |
| 3-trifluoromethyl phenoxy | 2-methylphenyl | 372 |
| 3-trifluoromethyl phenoxy | phenyl | 358 |
| 3-trifluoromethyl phenoxy | 4-ethylphenyl | 386 |
| 3-trifluoromethyl phenoxy | 2,3-difluorophenyl | 394 |
| 3-trifluoromethyl phenoxy | 2,6-dimethoxyphenyl | 418 |

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| 3-trifluoromethyl phenoxy | 3,4-difluorophenyl | 394 |
| 3-trifluoromethyl phenoxy | 2,5-difluorophenyl | 394 |
| 3-trifluoromethyl phenoxy | 4-ethoxyphenyl | 402 |
| 3-trifluoromethyl phenoxy | 2,4,6-trichlorophenyl | 462 |
| 3-trifluoromethyl phenoxy | 3-methylphenyl | 372 |
| 3-trifluoromethyl phenoxy | 2-fluoro-5- (trifluoromethyl)phenyl | 444 |
| 3-trifluoromethyl phenoxy | 3-methoxyphenyl | 388 |
| 3-trifluoromethyl phenoxy | 2-bromophenyl | 437 |
| 3-trifluoromethyl phenoxy | 4-bromophenyl | 437 |
| 3-trifluoromethyl phenoxy | 4-fluoro-3- (trifluoromethyl)phenyl | 444 |
| 3-trifluoromethyl phenoxy | 3-(trifluoromethoxy)phenyl | 442 |
| 3-trifluoromethyl phenoxy | 9-fluorenon-4-yl | 460 |
| 3-trifluoromethyl phenoxy | isoxazol-5-yl | 349 |
| 3-trifluoromethyl phenoxy | benzofuroxan-5-yl | 416 |
| 3-trifluoromethyl phenoxy | 2-chloropyrid-3-yl | 394 |
| 3-trifluoromethyl phenoxy | 2-(4- methylphenoxy)pyridin-3-yl | 465 |

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| 3-trifluoromethyl phenoxy | pyridin-4-yl | 359 |
| 3-trifluoromethyl phenoxy | anthraquinon-2-yl | 488 |
| 3-trifluoromethyl phenoxy | 2-iodophenyl | 484 |
| 3-trifluoromethyl phenoxy | 4-pentylphenyl | 428 |
| 3-trifluoromethyl phenoxy | 2-(4-chlorophenylthio) pyridin-3-yl | 502 |
| 3-trifluoromethyl phenoxy | 2,6-dimethylphenyl | 386 |
| 3-trifluoromethyl phenoxy | 2,5-dimethoxyphenyl | 418 |
| 3-trifluoromethyl phenoxy | 2,5-dichloropyridin-3-yl | 428 |
| 3-trifluoromethyl phenoxy | 2-chloro-6-methoxypyridin- 4-yl | 424 |
| 3-trifluoromethyl phenoxy | 2,3-dichloropyridin-5-yl | 428 |
| 3-trifluoromethyl phenoxy | 1-naphthyl | 422 |
| 3-trifluoromethyl phenoxy | 2,4-dimethoxyphenyl | 418 |
| 3-trifluoromethyl phenoxy | 3,5- bis(trifluoromethyl)phenyl | 494 |
| 3-trifluoromethyl phenoxy | 2-(4- chlorophenoxy)pyridin-3-yl | 486 |
| 3-trifluoromethyl phenoxy | pentafluorophenyl | 448 |
| 2,6-dichlorophenoxy | 3,4-dimethoxyphenyl | 419 |

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| 2,6-dichlorophenoxy | 2-(trifluoromethyl)phenyl | 427 |
| 2,6-dichlorophenoxy | 2,4-difluorophenyl | 395 |
| 2,6-dichlorophenoxy | 3-(trifluoromethyl)phenyl | 427 |
| 2,6-dichlorophenoxy | 2-naphthyl | 409 |
| 2,6-dichlorophenoxy | 2-methoxyphenyl | 389 |
| 2,6-dichlorophenoxy | 3,4,5-trimethylphenyl | 449 |
| 2,6-dichlorophenoxy | 3,4-dichlorophenyl | 428 |
| 2,6-dichlorophenoxy | 3-bromophenyl | 438 |
| 2,6-dichlorophenoxy | 3-pyridyl | 361 |
| 2,6-dichlorophenoxy | 2-ethoxynaphth-1-yl | 453 |
| 2,6-dichlorophenoxy | 2,3-dichlorophenyl | 428 |
| 2,6-dichlorophenoxy | 6-chloropyrid-3-yl | 395 |
| 2,6-dichlorophenoxy | 4-(trifluoromethoxy)phenyl | 443 |
| 2,6-dichlorophenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 445 |
| 2,6-dichlorophenoxy | 3-bromothienyl | 444 |
| 2,6-dichlorophenoxy | 2-acetoxyphenyl | 417 |
| 2,6-dichlorophenoxy | 5-methylisoxazol-3-yl | 364 |
| 2,6-dichlorophenoxy | 2-(phenylthio)pyrid-3-yl | 468 |
| 2,6-dichlorophenoxy | 2-(trifluoromethoxy)phenyl | 443 |
| 2,6-dichlorophenoxy | 1-phenyl-5-propylpyrazin-4-yl | 467 |
| 2,6-dichlorophenoxy | 2-ethoxyphenyl | 403 |
| 2,6-dichlorophenoxy | 3-chlorothien-2-yl | 400 |
| 2,6-dichlorophenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 419 |
| 2,6-dichlorophenoxy | 3,5-dichlorophenyl | 428 |
| 2,6-dichlorophenoxy | 2-(propylthio)pyridin-3-yl | 434 |

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| 2,6-dichlorophenoxy | 2-(ethylthio)pyridin-3-yl | 420 |
| 2,6-dichlorophenoxy | 3-bromopyridin-5-yl | 439 |
| 2,6-dichlorophenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 381 |
| 2,6-dichlorophenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 419 |
| 2,6-dichlorophenoxy | 3-chlorobenzo[b]thiophen-2-yl | 450 |
| 2,6-dichlorophenoxy | 4-chlorophenyl | 394 |
| 2,6-dichlorophenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 440 |
| 2,6-dichlorophenoxy | benzo[b]thiophen-2-yl | 415 |
| 2,6-dichlorophenoxy | 3,4-dimethylphenyl | 387 |
| 2,6-dichlorophenoxy | 2-(phenoxy)pyridin-3-yl | 452 |
| 2,6-dichlorophenoxy | 2-(methylthio)pyridin-3-yl | 406 |
| 2,6-dichlorophenoxy | 5-methyl-3-phenylisoxazol-4-yl | 440 |
| 2,6-dichlorophenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 463 |
| 2,6-dichlorophenoxy | 2-chloro-6-methylpyridin-4-yl | 409 |
| 2,6-dichlorophenoxy | 3,5-dimethylisoxazol-4-yl | 378 |
| 2,6-dichlorophenoxy | 1-naphthyl | 409 |
| 2,6-dichlorophenoxy | 2-fluorophenyl | 377 |
| 2,6-dichlorophenoxy | 4-propylphenyl | 401 |
| 2,6-dichlorophenoxy | 3-fluorophenyl | 377 |
| 2,6-dichlorophenoxy | 2,6-difluorophenyl | 395 |
| 2,6-dichlorophenoxy | 2-chlorophenyl | 394 |
| 2,6-dichlorophenoxy | 3-(chloromethyl)phenyl | 408 |

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| 2,6-dichlorophenoxy | 4-(2-(2-methyl)propyl)phenyl | 415 |
| 2,6-dichlorophenoxy | 3-chlorophenyl | 694 |
| 2,6-dichlorophenoxy | 3,5-dimethoxyphenyl | 419 |
| 2,6-dichlorophenoxy | 2,6-dichlorophenyl | 428 |
| 2,6-dichlorophenoxy | 2,4-dichlorophenyl | 428 |
| 2,6-dichlorophenoxy | 4-fluorophenyl | 377 |
| 2,6-dichlorophenoxy | 4-butylphenyl | 415 |
| 2,6-dichlorophenoxy | 2-methylphenyl | 373 |
| 2,6-dichlorophenoxy | phenyl | 359 |
| 2,6-dichlorophenoxy | 4-ethylphenyl | 387 |
| 2,6-dichlorophenoxy | 2,3-difluorophenyl | 395 |
| 2,6-dichlorophenoxy | 2,6-dimethoxyphenyl | 419 |
| 2,6-dichlorophenoxy | 3,4-difluorophenyl | 395 |
| 2,6-dichlorophenoxy | 2,5-difluorophenyl | 395 |
| 2,6-dichlorophenoxy | 4-ethoxyphenyl | 403 |
| 2,6-dichlorophenoxy | 2,4,6-trichlorophenyl | 463 |
| 2,6-dichlorophenoxy | 3-methylphenyl | 373 |
| 2,6-dichlorophenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 445 |
| 2,6-dichlorophenoxy | 3-methoxyphenyl | 389 |
| 2,6-dichlorophenoxy | 2-bromophenyl | 438 |
| 2,6-dichlorophenoxy | 4-bromophenyl | 438 |
| 2,6-dichlorophenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 445 |
| 2,6-dichlorophenoxy | 3-(trifluoromethoxy)phenyl | 443 |
| 2,6-dichlorophenoxy | 9-fluorenon-4-yl | 461 |
| 2,6-dichlorophenoxy | isoxazol-5-yl | 350 |

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| 2,6-dichlorophenoxy | benzofuroxan-5-yl | 417 |
| 2,6-dichlorophenoxy | 2-chloropyrid-3-yl | 395 |
| 2,6-dichlorophenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 466 |
| 2,6-dichlorophenoxy | pyridin-4-yl | 360 |
| 2,6-dichlorophenoxy | anthraquinon-2-yl | 489 |
| 2,6-dichlorophenoxy | 2-iodophenyl | 485 |
| 2,6-dichlorophenoxy | 4-pentylphenyl | 429 |
| 2,6-dichlorophenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 503 |
| 2,6-dichlorophenoxy | 2,6-dimethylphenyl | 387 |
| 2,6-dichlorophenoxy | 2,5-dimethoxyphenyl | 419 |
| 2,6-dichlorophenoxy | 2,5-dichloropyridin-3-yl | 429 |
| 2,6-dichlorophenoxy | 2-chloro-6-methoxypyridin-4-yl | 425 |
| 2,6-dichlorophenoxy | 2,3-dichloropyridin-5-yl | 429 |
| 2,6-dichlorophenoxy | 1-naphthyl | 413 |
| 2,6-dichlorophenoxy | 2,4-dimethoxyphenyl | 419 |
| 2,6-dichlorophenoxy | 3,5-bis(trifluoromethyl)phenyl | 495 |
| 2,6-dichlorophenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 487 |
| 2,6-dichlorophenoxy | pentafluorophenyl | 449 |
| 2,4-dichlorophenoxy | 3,4-dimethoxyphenyl | 419 |
| 2,4-dichlorophenoxy | 2-(trifluoromethyl)phenyl | 427 |
| 2,4-dichlorophenoxy | 2,4-difluorophenyl | 395 |
| 2,4-dichlorophenoxy | 3-(trifluoromethyl)phenyl | 427 |
| 2,4-dichlorophenoxy | 2-naphthyl | 409 |
| 2,4-dichlorophenoxy | 2-methoxyphenyl | 389 |

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| 2,4-dichlorophenoxy | 3,4,5-trimethylphenyl | 449 |
| 2,4-dichlorophenoxy | 3,4-dichlorophenyl | 428 |
| 2,4-dichlorophenoxy | 3-bromophenyl | 438 |
| 2,4-dichlorophenoxy | 3-pyridyl | 361 |
| 2,4-dichlorophenoxy | 2-ethoxynaphth-1-yl | 453 |
| 2,4-dichlorophenoxy | 2,3-dichlorophenyl | 428 |
| 2,4-dichlorophenoxy | 6-chloropyrid-3-yl | 395 |
| 2,4-dichlorophenoxy | 4-(trifluoromethoxy)phenyl | 443 |
| 2,4-dichlorophenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 445 |
| 2,4-dichlorophenoxy | 3-bromothieryl | 444 |
| 2,4-dichlorophenoxy | 2-acetoxyphenyl | 417 |
| 2,4-dichlorophenoxy | 5-methylisoxazol-3-yl | 364 |
| 2,4-dichlorophenoxy | 2-(phenylthio)pyrid-3-yl | 468 |
| 2,4-dichlorophenoxy | 2-(trifluoromethoxy)phenyl | 443 |
| 2,4-dichlorophenoxy | 1-phenyl-5-propylpyrazin-4-yl | 467 |
| 2,4-dichlorophenoxy | 2-ethoxyphenyl | 403 |
| 2,4-dichlorophenoxy | 3-chlorothieryl-2-yl | 400 |
| 2,4-dichlorophenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 419 |
| 2,4-dichlorophenoxy | 3,5-dichlorophenyl | 428 |
| 2,4-dichlorophenoxy | 2-(propylthio)pyridin-3-yl | 434 |
| 2,4-dichlorophenoxy | 2-(ethylthio)pyridin-3-yl | 420 |
| 2,4-dichlorophenoxy | 3-bromopyridin-5-yl | 439 |
| 2,4-dichlorophenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 381 |
| 2,4-dichlorophenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 419 |

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| 2,4-dichlorophenoxy | 3-chlorobenzo[b]thiophen-2-yl | 450 |
| 2,4-dichlorophenoxy | 4-chlorophenyl | 394 |
| 2,4-dichlorophenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 440 |
| 2,4-dichlorophenoxy | benzo[b]thiophen-2-yl | 415 |
| 2,4-dichlorophenoxy | 3,4-dimethylphenyl | 387 |
| 2,4-dichlorophenoxy | 2-(phenoxy)pyridin-3-yl | 452 |
| 2,4-dichlorophenoxy | 2-(methylthio)pyridin-3-yl | 406 |
| 2,4-dichlorophenoxy | 5-methyl-3-phenylisoxazol-4-yl | 440 |
| 2,4-dichlorophenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 463 |
| 2,4-dichlorophenoxy | 2-chloro-6-methylpyridin-4-yl | 409 |
| 2,4-dichlorophenoxy | 3,5-dimethylisoxazol-4-yl | 378 |
| 2,4-dichlorophenoxy | 1-naphthyl | 409 |
| 2,4-dichlorophenoxy | 2-fluorophenyl | 377 |
| 2,4-dichlorophenoxy | 4-propylphenyl | 401 |
| 2,4-dichlorophenoxy | 3-fluorophenyl | 377 |
| 2,4-dichlorophenoxy | 2,6-difluorophenyl | 395 |
| 2,4-dichlorophenoxy | 2-chlorophenyl | 394 |
| 2,4-dichlorophenoxy | 3-(chloromethyl)phenyl | 408 |
| 2,4-dichlorophenoxy | 4-(2-(2-methyl)propyl)phenyl | 415 |
| 2,4-dichlorophenoxy | 3-chlorophenyl | 694 |
| 2,4-dichlorophenoxy | 3,5-dimethoxyphenyl | 419 |
| 2,4-dichlorophenoxy | 2,6-dichlorophenyl | 428 |

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| 2,4-dichlorophenoxy | 2,4-dichlorophenyl | 428 |
| 2,4-dichlorophenoxy | 4-fluorophenyl | 377 |
| 2,4-dichlorophenoxy | 4-butylphenyl | 415 |
| 2,4-dichlorophenoxy | 2-methylphenyl | 373 |
| 2,4-dichlorophenoxy | phenyl | 359 |
| 2,4-dichlorophenoxy | 4-ethylphenyl | 387 |
| 2,4-dichlorophenoxy | 2,3-difluorophenyl | 395 |
| 2,4-dichlorophenoxy | 2,6-dimethoxyphenyl | 419 |
| 2,4-dichlorophenoxy | 3,4-difluorophenyl | 395 |
| 2,4-dichlorophenoxy | 2,5-difluorophenyl | 395 |
| 2,4-dichlorophenoxy | 4-ethoxyphenyl | 403 |
| 2,4-dichlorophenoxy | 2,4,6-trichlorophenyl | 463 |
| 2,4-dichlorophenoxy | 3-methylphenyl | 373 |
| 2,4-dichlorophenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 445 |
| 2,4-dichlorophenoxy | 3-methoxyphenyl | 389 |
| 2,4-dichlorophenoxy | 2-bromophenyl | 438 |
| 2,4-dichlorophenoxy | 4-bromophenyl | 438 |
| 2,4-dichlorophenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 445 |
| 2,4-dichlorophenoxy | 3-(trifluoromethoxy)phenyl | 443 |
| 2,4-dichlorophenoxy | 9-fluorenon-4-yl | 461 |
| 2,4-dichlorophenoxy | isoxazol-5-yl | 350 |
| 2,4-dichlorophenoxy | benzofuroxan-5-yl | 417 |
| 2,4-dichlorophenoxy | 2-chloropyrid-3-yl | 395 |
| 2,4-dichlorophenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 466 |
| 2,4-dichlorophenoxy | pyridin-4-yl | 360 |

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| 2,4-dichlorophenoxy | anthraquinon-2-yl | 489 |
| 2,4-dichlorophenoxy | 2-iodophenyl | 485 |
| 2,4-dichlorophenoxy | 4-pentylphenyl | 429 |
| 2,4-dichlorophenoxy | 2-(4-chlorophenylthio) pyridin-3-yl | 503 |
| 2,4-dichlorophenoxy | 2,6-dimethylphenyl | 387 |
| 2,4-dichlorophenoxy | 2,5-dimethoxyphenyl | 419 |
| 2,4-dichlorophenoxy | 2,5-dichloropyridin-3-yl | 429 |
| 2,4-dichlorophenoxy | 2-chloro-6-methoxypyridin- 4-yl | 425 |
| 2,4-dichlorophenoxy | 2,3-dichloropyridin-5-yl | 429 |
| 2,4-dichlorophenoxy | 1-naphthyl | 413 |
| 2,4-dichlorophenoxy | 2,4-dimethoxyphenyl | 419 |
| 2,4-dichlorophenoxy | 3,5- bis(trifluoromethyl)phenyl | 495 |
| 2,4-dichlorophenoxy | 2-(4- chlorophenoxy)pyridin-3-yl | 487 |
| 2,4-dichlorophenoxy | pentafluorophenyl | 449 |
| 4-chloro-3- methylphenoxy | 3,4-dimethoxyphenyl | 319 |
| 4-chloro-3- methylphenoxy | 2-(trifluoromethyl)phenyl | 407 |
| 4-chloro-3- methylphenoxy | 2,4-difluorophenyl | 375 |
| 4-chloro-3- methylphenoxy | 3-(trifluoromethyl)phenyl | 407 |
| 4-chloro-3- methylphenoxy | 2-naphthyl | 389 |
| 4-chloro-3- methylphenoxy | 2-methoxyphenyl | 369 |

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| 4-chloro-3-methylphenoxy | 3,4,5-trimethylphenyl | 429 |
| 4-chloro-3-methylphenoxy | 3,4-dichlorophenyl | 408 |
| 4-chloro-3-methylphenoxy | 3-bromophenyl | 418 |
| 4-chloro-3-methylphenoxy | 3-pyridyl | 340 |
| 4-chloro-3-methylphenoxy | 2-ethoxynaphth-1-yl | 433 |
| 4-chloro-3-methylphenoxy | 2,3-dichlorophenyl | 408 |
| 4-chloro-3-methylphenoxy | 6-chloropyrid-3-yl | 374 |
| 4-chloro-3-methylphenoxy | 4-(trifluoromethoxy)phenyl | 423 |
| 4-chloro-3-methylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 425 |
| 4-chloro-3-methylphenoxy | 3-bromothienyl | 424 |
| 4-chloro-3-methylphenoxy | 2-acetoxyphenyl | 397 |
| 4-chloro-3-methylphenoxy | 5-methylisoxazol-3-yl | 344 |
| 4-chloro-3-methylphenoxy | 2-(phenylthio)pyrid-3-yl | 448 |
| 4-chloro-3-methylphenoxy | 2-(trifluoromethoxy)phenyl | 423 |
| 4-chloro-3-methylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 447 |
| 4-chloro-3-methylphenoxy | 2-ethoxyphenyl | 383 |

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| 4-chloro-3-methylphenoxy | 3-chlorothien-2-yl | 379 |
| 4-chloro-3-methylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 399 |
| 4-chloro-3-methylphenoxy | 3,5-dichlorophenyl | 408 |
| 4-chloro-3-methylphenoxy | 2-(propylthio)pyridin-3-yl | 414 |
| 4-chloro-3-methylphenoxy | 2-(ethylthio)pyridin-3-yl | 400 |
| 4-chloro-3-methylphenoxy | 3-bromopyridin-5-yl | 419 |
| 4-chloro-3-methylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 361 |
| 4-chloro-3-methylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 399 |
| 4-chloro-3-methylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 429 |
| 4-chloro-3-methylphenoxy | 4-chlorophenyl | 373 |
| 4-chloro-3-methylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 420 |
| 4-chloro-3-methylphenoxy | benzo[b]thiophen-2-yl | 395 |
| 4-chloro-3-methylphenoxy | 3,4-dimethylphenyl | 367 |
| 4-chloro-3-methylphenoxy | 2-(phenoxy)pyridin-3-yl | 432 |
| 4-chloro-3-methylphenoxy | 2-(methylthio)pyridin-3-yl | 386 |
| 4-chloro-3-methylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 420 |

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| 4-chloro-3-methylphenoxy | 4-chloro-1,3-dimethyl pyrazolo[3,4-b]pyridin-3-yl | 442 |
| 4-chloro-3-methylphenoxy | 2-chloro-6-methylpyridin-4-yl | 388 |
| 4-chloro-3-methylphenoxy | 3,5-dimethylisoxazol-4-yl | 358 |
| 4-chloro-3-methylphenoxy | 1-naphthyl | 389 |
| 4-chloro-3-methylphenoxy | 2-fluorophenyl | 357 |
| 4-chloro-3-methylphenoxy | 4-propylphenyl | 381 |
| 4-chloro-3-methylphenoxy | 4-(trifluoromethyl)phenyl | 407 |
| 4-chloro-3-methylphenoxy | 3-fluorophenyl | 357 |
| 4-chloro-3-methylphenoxy | 2,6-difluorophenyl | 375 |
| 4-chloro-3-methylphenoxy | 2-chlorophenyl | 373 |
| 4-chloro-3-methylphenoxy | 3-(chloromethyl)phenyl | 387 |
| 4-chloro-3-methylphenoxy | 4-(2-(2-methyl)propyl)phenyl | 395 |
| 4-chloro-3-methylphenoxy | 3-chlorophenyl | 373 |
| 4-chloro-3-methylphenoxy | 3,5-dimethoxyphenyl | 399 |
| 4-chloro-3-methylphenoxy | 2,6-dichlorophenyl | 408 |

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| 4-chloro-3-methylphenoxy | 2,4-dichlorophenyl | 408 |
| 4-chloro-3-methylphenoxy | 4-fluorophenyl | 357 |
| 4-chloro-3-methylphenoxy | 4-butylphenyl | 395 |
| 4-chloro-3-methylphenoxy | 2-methylphenyl | 353 |
| 4-chloro-3-methylphenoxy | phenyl | 339 |
| 4-chloro-3-methylphenoxy | 4-ethylphenyl | 367 |
| 4-chloro-3-methylphenoxy | 2,3-difluorophenyl | 375 |
| 4-chloro-3-methylphenoxy | 2,6-dimethoxyphenyl | 399 |
| 4-chloro-3-methylphenoxy | 3,4-difluorophenyl | 375 |
| 4-chloro-3-methylphenoxy | 2,5-difluorophenyl | 375 |
| 4-chloro-3-methylphenoxy | 4-ethoxyphenyl | 383 |
| 4-chloro-3-methylphenoxy | 2,4,6-trichlorophenyl | 442 |
| 4-chloro-3-methylphenoxy | 3-methylphenyl | 353 |
| 4-chloro-3-methylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 425 |
| 4-chloro-3-methylphenoxy | 3-methoxyphenyl | 369 |

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| 4-chloro-3-methylphenoxy | 2-bromophenyl | 418 |
| 4-chloro-3-methylphenoxy | 4-bromophenyl | 418 |
| 4-chloro-3-methylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 425 |
| 4-chloro-3-methylphenoxy | 3-(trifluoromethoxy)phenyl | 423 |
| 4-chloro-3-methylphenoxy | 9-fluorenon-4-yl | 441 |
| 4-chloro-3-methylphenoxy | isoxazol-5-yl | 330 |
| 4-chloro-3-methylphenoxy | benzofuroxan-5-yl | 397 |
| 4-chloro-3-methylphenoxy | 2-chloropyrid-3-yl | 374 |
| 4-chloro-3-methylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 446 |
| 4-chloro-3-methylphenoxy | pyridin-4-yl | 340 |
| 4-chloro-3-methylphenoxy | anthraquinon-2-yl | 469 |
| 4-chloro-3-methylphenoxy | 2-iodophenyl | 465 |
| 4-chloro-3-methylphenoxy | 4-pentylphenyl | 409 |
| 4-chloro-3-methylphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 482 |
| 4-chloro-3-methylphenoxy | 2,6-dimethylphenyl | 367 |

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| 4-chloro-3-methylphenoxy | 2,5-dimethoxyphenyl | 399 |
| 4-chloro-3-methylphenoxy | 2,5-dichloropyridin-3-yl | 409 |
| 4-chloro-3-methylphenoxy | 2-chloro-6-methoxypyridin-4-yl | 404 |
| 4-chloro-3-methylphenoxy | 2,3-dichloropyridin-5-yl | 409 |
| 4-chloro-3-methylphenoxy | 1-naphthyl | 403 |
| 4-chloro-3-methylphenoxy | 2,4-dimethoxyphenyl | 399 |
| 4-chloro-3-methylphenoxy | 3,5-bis(trifluoromethyl)phenyl | 475 |
| 4-chloro-3-methylphenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 466 |
| 4-chloro-3-methylphenoxy | pentafluorophenyl | 429 |
| 4-chloro-2-cyclohexylphenoxy | 3,4-dimethoxyphenyl | 467 |
| 4-chloro-2-cyclohexylphenoxy | 2-(trifluoromethyl)phenyl | 475 |
| 4-chloro-2-cyclohexylphenoxy | 2,4-difluorophenyl | 443 |
| 4-chloro-2-cyclohexylphenoxy | 3-(trifluoromethyl)phenyl | 475 |
| 4-chloro-2-cyclohexylphenoxy | 2-naphthyl | 457 |
| 4-chloro-2-cyclohexylphenoxy | 2-methoxyphenyl | 437 |

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| 4-chloro-2-cyclohexylphenoxy | 3,4,5-trimethylphenyl | 497 |
| 4-chloro-2-cyclohexylphenoxy | 3,4-dichlorophenyl | 176 |
| 4-chloro-2-cyclohexylphenoxy | 3-bromophenyl | 486 |
| 4-chloro-2-cyclohexylphenoxy | 3-pyridyl | 408 |
| 4-chloro-2-cyclohexylphenoxy | 2-ethoxynaphth-1-yl | 501 |
| 4-chloro-2-cyclohexylphenoxy | 2,3-dichlorophenyl | 476 |
| 4-chloro-2-cyclohexylphenoxy | 6-chloropyrid-3-yl | 442 |
| 4-chloro-2-cyclohexylphenoxy | 4-(trifluoromethoxy)phenyl | 491 |
| 4-chloro-2-cyclohexylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 493 |
| 4-chloro-2-cyclohexylphenoxy | 3-bromothieryl | 492 |
| 4-chloro-2-cyclohexylphenoxy | 2-acetoxyphenyl | 465 |
| 4-chloro-2-cyclohexylphenoxy | 5-methylisoxazol-3-yl | 412 |
| 4-chloro-2-cyclohexylphenoxy | 2-(phenylthio)pyrid-3-yl | 516 |
| 4-chloro-2-cyclohexylphenoxy | 2-(trifluoromethoxy)phenyl | 491 |
| 4-chloro-2-cyclohexylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 515 |

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| 4-chloro-2-cyclohexylphenoxy | 2-ethoxyphenyl | 451 |
| 4-chloro-2-cyclohexylphenoxy | 3-chlorothien-2-yl | 447 |
| 4-chloro-2-cyclohexylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 467 |
| 4-chloro-2-cyclohexylphenoxy | 3,5-dichlorophenyl | 476 |
| 4-chloro-2-cyclohexylphenoxy | 2-(propylthio)pyridin-3-yl | 482 |
| 4-chloro-2-cyclohexylphenoxy | 2-(ethylthio)pyridin-3-yl | 468 |
| 4-chloro-2-cyclohexylphenoxy | 3-bromopyridin-5-yl | 487 |
| 4-chloro-2-cyclohexylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 429 |
| 4-chloro-2-cyclohexylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 467 |
| 4-chloro-2-cyclohexylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 497 |
| 4-chloro-2-cyclohexylphenoxy | 4-chlorophenyl | 441 |
| 4-chloro-2-cyclohexylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 488 |
| 4-chloro-2-cyclohexylphenoxy | benzo[b]thiophen-2-yl | 463 |
| 4-chloro-2-cyclohexylphenoxy | 3,4-dimethylphenyl | 435 |
| 4-chloro-2-cyclohexylphenoxy | 2-(phenoxy)pyridin-3-yl | 500 |

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| 4-chloro-2-cyclohexylphenoxy | 2-(methylthio)pyridin-3-yl | 454 |
| 4-chloro-2-cyclohexylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 488 |
| 4-chloro-2-cyclohexylphenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 510 |
| 4-chloro-2-cyclohexylphenoxy | 2-chloro-6-methylpyridin-4-yl | 456 |
| 4-chloro-2-cyclohexylphenoxy | 3,5-dimethylisoxazol-4-yl | 426 |
| 4-chloro-2-cyclohexylphenoxy | 1-naphthyl | 457 |
| 4-chloro-2-cyclohexylphenoxy | 2-fluorophenyl | 425 |
| 4-chloro-2-cyclohexylphenoxy | 4-propylphenyl | 449 |
| 4-chloro-2-cyclohexylphenoxy | 3-fluorophenyl | 425 |
| 4-chloro-2-cyclohexylphenoxy | 2,6-difluorophenyl | 443 |
| 4-chloro-2-cyclohexylphenoxy | 2-chlorophenyl | 441 |
| 4-chloro-2-cyclohexylphenoxy | 3-(chloromethyl)phenyl | 455 |
| 4-chloro-2-cyclohexylphenoxy | 4-(2-(2-methyl)propyl)phenyl | 463 |
| 4-chloro-2-cyclohexylphenoxy | 3-chlorophenyl | 441 |
| 4-chloro-2-cyclohexylphenoxy | 3,5-dimethoxyphenyl | 467 |

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| 4-chloro-2-cyclohexylphenoxy | 2,6-dichlorophenyl | 476 |
| 4-chloro-2-cyclohexylphenoxy | 2,4-dichlorophenyl | 476 |
| 4-chloro-2-cyclohexylphenoxy | 4-fluorophenyl | 425 |
| 4-chloro-2-cyclohexylphenoxy | 4-butylphenyl | 463 |
| 4-chloro-2-cyclohexylphenoxy | 2-methylphenyl | 421 |
| 4-chloro-2-cyclohexylphenoxy | phenyl | 407 |
| 4-chloro-2-cyclohexylphenoxy | 4-ethylphenyl | 435 |
| 4-chloro-2-cyclohexylphenoxy | 2,3-difluorophenyl | 443 |
| 4-chloro-2-cyclohexylphenoxy | 2,6-dimethoxyphenyl | 467 |
| 4-chloro-2-cyclohexylphenoxy | 3,4-difluorophenyl | 443 |
| 4-chloro-2-cyclohexylphenoxy | 2,5-difluorophenyl | 443 |
| 4-chloro-2-cyclohexylphenoxy | 4-ethoxyphenyl | 451 |
| 4-chloro-2-cyclohexylphenoxy | 2,4,6-trichlorophenyl | 510 |
| 4-chloro-2-cyclohexylphenoxy | 3-methylphenyl | 421 |
| 4-chloro-2-cyclohexylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 493 |
| 4-chloro-2-cyclohexylphenoxy | 3-methoxyphenyl | 437 |

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| 4-chloro-2-cyclohexylphenoxy | 2-bromophenyl | 486 |
| 4-chloro-2-cyclohexylphenoxy | 4-bromophenyl | 486 |
| 4-chloro-2-cyclohexylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 493 |
| 4-chloro-2-cyclohexylphenoxy | 3-(trifluoromethoxy)phenyl | 491 |
| 4-chloro-2-cyclohexylphenoxy | 9-fluorenon-4-yl | 503 |
| 4-chloro-2-cyclohexylphenoxy | isoxazol-5-yl | 398 |
| 4-chloro-2-cyclohexylphenoxy | benzofuroxan-5-yl | 465 |
| 4-chloro-2-cyclohexylphenoxy | 2-chloropyrid-3-yl | 442 |
| 4-chloro-2-cyclohexylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 514 |
| 4-chloro-2-cyclohexylphenoxy | pyridin-4-yl | 408 |
| 4-chloro-2-cyclohexylphenoxy | anthraquinon-2-yl | 537 |
| 4-chloro-2-cyclohexylphenoxy | 2-iodophenyl | 533 |
| 4-chloro-2-cyclohexylphenoxy | 4-pentylphenyl | 477 |
| 4-chloro-2-cyclohexylphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 550 |
| 4-chloro-2-cyclohexylphenoxy | 2,6-dimethylphenyl | 435 |
| 4-chloro-2-cyclohexylphenoxy | 2,5-dimethoxyphenyl | 467 |

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| 4-chloro-2-cyclohexylphenoxy | 2,5-dichloropyridin-3-yl | 477 |
| 4-chloro-2-cyclohexylphenoxy | 2-chloro-6-methoxypyridin-4-yl | 472 |
| 4-chloro-2-cyclohexylphenoxy | 2,3-dichloropyridin-5-yl | 477 |
| 4-chloro-2-cyclohexylphenoxy | 1-naphthyl | 471 |
| 4-chloro-2-cyclohexylphenoxy | 2,4-dimethoxyphenyl | 467 |
| 4-chloro-2-cyclohexylphenoxy | 3,5-bis(trifluoromethyl)phenyl | 546 |
| 4-chloro-2-cyclohexylphenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 534 |
| 4-chloro-2-cyclohexylphenoxy | pentafluorophenyl | 497 |
| 4-chloro-3,5-dimethylphenoxy | 3,4-dimethoxyphenyl | 413 |
| 4-chloro-3,5-dimethylphenoxy | 2-(trifluoromethyl)phenyl | 421 |
| 4-chloro-3,5-dimethylphenoxy | 2,4-difluorophenyl | 389 |
| 4-chloro-3,5-dimethylphenoxy | 3-(trifluoromethyl)phenyl | 421 |
| 4-chloro-3,5-dimethylphenoxy | 2-naphthyl | 403 |
| 4-chloro-3,5-dimethylphenoxy | 2-methoxyphenyl | 484 |
| 4-chloro-3,5-dimethylphenoxy | 3,4,5-trimethylphenyl | 443 |
| 4-chloro-3,5-dimethylphenoxy | 3,4-dichlorophenyl | 422 |

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| 4-chloro-3,5-dimethylphenoxy | 3-bromophenyl | 432 |
| 4-chloro-3,5-dimethylphenoxy | 3-pyridyl | 354 |
| 4-chloro-3,5-dimethylphenoxy | 2-ethoxynaphth-1-yl | 447 |
| 4-chloro-3,5-dimethylphenoxy | 2,3-dichlorophenyl | 422 |
| 4-chloro-3,5-dimethylphenoxy | 6-chloropyrid-3-yl | 388 |
| 4-chloro-3,5-dimethylphenoxy | 4-(trifluoromethoxy)phenyl | 437 |
| 4-chloro-3,5-dimethylphenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 439 |
| 4-chloro-3,5-dimethylphenoxy | 3-bromothienyl | 438 |
| 4-chloro-3,5-dimethylphenoxy | 2-acetoxyphenyl | 411 |
| 4-chloro-3,5-dimethylphenoxy | 5-methylisoxazol-3-yl | 358 |
| 4-chloro-3,5-dimethylphenoxy | 2-(phenylthio)pyrid-3-yl | 462 |
| 4-chloro-3,5-dimethylphenoxy | 2-(trifluoromethoxy)phenyl | 437 |
| 4-chloro-3,5-dimethylphenoxy | 1-phenyl-5-propylpyrazin-4-yl | 461 |
| 4-chloro-3,5-dimethylphenoxy | 2-ethoxyphenyl | 397 |
| 4-chloro-3,5-dimethylphenoxy | 3-chlorothien-2-yl | 393 |
| 4-chloro-3,5-dimethylphenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 413 |

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| 4-chloro-3,5-dimethylphenoxy | 3,5-dichlorophenyl | 422 |
| 4-chloro-3,5-dimethylphenoxy | 2-(propylthio)pyridin-3-yl | 428 |
| 4-chloro-3,5-dimethylphenoxy | 2-(ethylthio)pyridin-3-yl | 414 |
| 4-chloro-3,5-dimethylphenoxy | 3-bromopyridin-5-yl | 433 |
| 4-chloro-3,5-dimethylphenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 375 |
| 4-chloro-3,5-dimethylphenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 413 |
| 4-chloro-3,5-dimethylphenoxy | 3-chlorobenzo[b]thiophen-2-yl | 443 |
| 4-chloro-3,5-dimethylphenoxy | 4-chlorophenyl | 387 |
| 4-chloro-3,5-dimethylphenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 434 |
| 4-chloro-3,5-dimethylphenoxy | benzo[b]thiophen-2-yl | 409 |
| 4-chloro-3,5-dimethylphenoxy | 3,4-dimethylphenyl | 381 |
| 4-chloro-3,5-dimethylphenoxy | 2-(phenoxy)pyridin-3-yl | 446 |
| 4-chloro-3,5-dimethylphenoxy | 2-(methylthio)pyridin-3-yl | 400 |
| 4-chloro-3,5-dimethylphenoxy | 5-methyl-3-phenylisoxazol-4-yl | 434 |
| 4-chloro-3,5-dimethylphenoxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 456 |

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| 4-chloro-3,5-dimethylphenoxy | 2-chloro-6-methylpyridin-4-yl | 402 |
| 4-chloro-3,5-dimethylphenoxy | 3,5-dimethylisoxazol-4-yl | 372 |
| 4-chloro-3,5-dimethylphenoxy | 1-naphthyl | 403 |
| 4-chloro-3,5-dimethylphenoxy | 2-fluorophenyl | 371 |
| 4-chloro-3,5-dimethylphenoxy | 4-propylphenyl | 395 |
| 4-chloro-3,5-dimethylphenoxy | 3-fluorophenyl | 371 |
| 4-chloro-3,5-dimethylphenoxy | 2,6-difluorophenyl | 389 |
| 4-chloro-3,5-dimethylphenoxy | 2-chlorophenyl | 387 |
| 4-chloro-3,5-dimethylphenoxy | 3-(chloromethyl)phenyl | 401 |
| 4-chloro-3,5-dimethylphenoxy | 4-(2-(2-methyl)propyl)phenyl | 409 |
| 4-chloro-3,5-dimethylphenoxy | 3-chlorophenyl | 387 |
| 4-chloro-3,5-dimethylphenoxy | 3,5-dimethoxyphenyl | 413 |
| 4-chloro-3,5-dimethylphenoxy | 2,6-dichlorophenyl | 422 |
| 4-chloro-3,5-dimethylphenoxy | 2,4-dichlorophenyl | 422 |
| 4-chloro-3,5-dimethylphenoxy | 4-fluorophenyl | 371 |

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| 4-chloro-3,5-dimethylphenoxy | 4-butylphenyl | 409 |
| 4-chloro-3,5-dimethylphenoxy | 2-methylphenyl | 367 |
| 4-chloro-3,5-dimethylphenoxy | phenyl | 353 |
| 4-chloro-3,5-dimethylphenoxy | 4-ethylphenyl | 381 |
| 4-chloro-3,5-dimethylphenoxy | 2,3-difluorophenyl | 389 |
| 4-chloro-3,5-dimethylphenoxy | 2,6-dimethoxyphenyl | 413 |
| 4-chloro-3,5-dimethylphenoxy | 3,4-difluorophenyl | 389 |
| 4-chloro-3,5-dimethylphenoxy | 2,5-difluorophenyl | 389 |
| 4-chloro-3,5-dimethylphenoxy | 4-ethoxyphenyl | 397 |
| 4-chloro-3,5-dimethylphenoxy | 2,4,6-trichlorophenyl | 456 |
| 4-chloro-3,5-dimethylphenoxy | 3-methylphenyl | 367 |
| 4-chloro-3,5-dimethylphenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 439 |
| 4-chloro-3,5-dimethylphenoxy | 3-methoxyphenyl | 383 |
| 4-chloro-3,5-dimethylphenoxy | 2-bromophenyl | 432 |
| 4-chloro-3,5-dimethylphenoxy | 4-bromophenyl | 432 |

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| 4-chloro-3,5-dimethylphenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 439 |
| 4-chloro-3,5-dimethylphenoxy | 3-(trifluoromethoxy)phenyl | 437 |
| 4-chloro-3,5-dimethylphenoxy | 9-fluorenon-4-yl | 455 |
| 4-chloro-3,5-dimethylphenoxy | isoxazol-5-yl | 344 |
| 4-chloro-3,5-dimethylphenoxy | benzofuroxan-5-yl | 411 |
| 4-chloro-3,5-dimethylphenoxy | 2-chloropyrid-3-yl | 388 |
| 4-chloro-3,5-dimethylphenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 460 |
| 4-chloro-3,5-dimethylphenoxy | pyridin-4-yl | 354 |
| 4-chloro-3,5-dimethylphenoxy | anthraquinon-2-yl | 483 |
| 4-chloro-3,5-dimethylphenoxy | 2-iodophenyl | 479 |
| 4-chloro-3,5-dimethylphenoxy | 4-pentylphenyl | 423 |
| 4-chloro-3,5-dimethylphenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 496 |
| 4-chloro-3,5-dimethylphenoxy | 2,6-dimethylphenyl | 381 |
| 4-chloro-3,5-dimethylphenoxy | 2,5-dimethoxyphenyl | 413 |
| 4-chloro-3,5-dimethylphenoxy | 2,5-dichloropyridin-3-yl | 423 |

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| 4-chloro-3,5-dimethylphenoxy | 2-chloro-6-methoxypyridin-4-yl | 418 |
| 4-chloro-3,5-dimethylphenoxy | 2,3-dichloropyridin-5-yl | 423 |
| 4-chloro-3,5-dimethylphenoxy | 1-naphthyl | 417 |
| 4-chloro-3,5-dimethylphenoxy | 2,4-dimethoxyphenyl | 413 |
| 4-chloro-3,5-dimethylphenoxy | 3,5-bis(trifluoromethyl)phenyl | 489 |
| 4-chloro-3,5-dimethylphenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 480 |
| 4-chloro-3,5-dimethylphenoxy | pentafluorophenyl | 443 |
| pyrid-3-yloxy | 3,4-dimethoxyphenyl | 351 |
| pyrid-3-yloxy | 2-(trifluoromethyl)phenyl | 359 |
| pyrid-3-yloxy | 2,4-difluorophenyl | 327 |
| pyrid-3-yloxy | 3-(trifluoromethyl)phenyl | 359 |
| pyrid-3-yloxy | 2-naphthyl | 341 |
| pyrid-3-yloxy | 2-methoxyphenyl | 321 |
| pyrid-3-yloxy | 3,4,5-trimethylphenyl | 381 |
| pyrid-3-yloxy | 3,4-dichlorophenyl | 360 |
| pyrid-3-yloxy | 3-bromophenyl | 370 |
| pyrid-3-yloxy | 3-pyridyl | 292 |
| pyrid-3-yloxy | 2-ethoxynaphth-1-yl | 385 |
| pyrid-3-yloxy | 2,3-dichlorophenyl | 360 |
| pyrid-3-yloxy | 6-chloropyrid-3-yl | 327 |
| pyrid-3-yloxy | 4-(trifluoromethoxy)phenyl | 375 |

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| pyrid-3-yloxy | 2-fluoro-4-(trifluoromethyl)phenyl | 377 |
| pyrid-3-yloxy | 3-bromothienyl | 376 |
| pyrid-3-yloxy | 2-acetoxyphenyl | 349 |
| pyrid-3-yloxy | 5-methylisoxazol-3-yl | 296 |
| pyrid-3-yloxy | 2-(phenylthio)pyrid-3-yl | 400 |
| pyrid-3-yloxy | 2-(trifluoromethoxy)phenyl | 375 |
| pyrid-3-yloxy | 1-phenyl-5-propylpyrazin-4-yl | 399 |
| pyrid-3-yloxy | 2-ethoxyphenyl | 335 |
| pyrid-3-yloxy | 3-chlorothien-2-yl | 332 |
| pyrid-3-yloxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 351 |
| pyrid-3-yloxy | 3,5-dichlorophenyl | 360 |
| pyrid-3-yloxy | 2-(propylthio)pyridin-3-yl | 366 |
| pyrid-3-yloxy | 2-(ethylthio)pyridin-3-yl | 352 |
| pyrid-3-yloxy | 3-bromopyridin-5-yl | 371 |
| pyrid-3-yloxy | 4-methyl-1,2,3-thiadiazol-5-yl | 313 |
| pyrid-3-yloxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 351 |
| pyrid-3-yloxy | 3-chlorobenzo[b]thiophen-2-yl | 382 |
| pyrid-3-yloxy | 4-chlorophenyl | 326 |
| pyrid-3-yloxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 372 |
| pyrid-3-yloxy | benzo[b]thiophen-2-yl | 347 |
| pyrid-3-yloxy | 3,4-dimethylphenyl | 319 |
| pyrid-3-yloxy | 2-(phenoxy)pyridin-3-yl | 384 |

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| pyrid-3-yloxy | 2-(methylthio)pyridin-3-yl | 338 |
| pyrid-3-yloxy | 5-methyl-3-phenylisoxazol-4-yl | 372 |
| pyrid-3-yloxy | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 395 |
| pyrid-3-yloxy | 2-chloro-6-methylpyridin-4-yl | 341 |
| pyrid-3-yloxy | 3,5-dimethylisoxazol-4-yl | 310 |
| pyrid-3-yloxy | 1-naphthyl | 341 |
| pyrid-3-yloxy | 2-fluorophenyl | 309 |
| pyrid-3-yloxy | 4-propylphenyl | 333 |
| pyrid-3-yloxy | 3-fluorophenyl | 309 |
| pyrid-3-yloxy | 2,6-difluorophenyl | 327 |
| pyrid-3-yloxy | 2-chlorophenyl | 326 |
| pyrid-3-yloxy | 3-(chloromethyl)phenyl | 340 |
| pyrid-3-yloxy | 4-(2-(2-methyl)propyl)phenyl | 347 |
| pyrid-3-yloxy | 3-chlorophenyl | 326 |
| pyrid-3-yloxy | 3,5-dimethoxyphenyl | 351 |
| pyrid-3-yloxy | 2,6-dichlorophenyl | 360 |
| pyrid-3-yloxy | 2,4-dichlorophenyl | 360 |
| pyrid-3-yloxy | 4-fluorophenyl | 309 |
| pyrid-3-yloxy | 4-butylphenyl | 347 |
| pyrid-3-yloxy | 2-methylphenyl | 305 |
| pyrid-3-yloxy | phenyl | 291 |
| pyrid-3-yloxy | 4-ethylphenyl | 319 |
| pyrid-3-yloxy | 2,3-difluorophenyl | 327 |
| pyrid-3-yloxy | 2,6-dimethoxyphenyl | 351 |

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| pyrid-3-yloxy | 3,4-difluorophenyl | 327 |
| pyrid-3-yloxy | 2,5-difluorophenyl | 327 |
| pyrid-3-yloxy | 4-ethoxyphenyl | 335 |
| pyrid-3-yloxy | 2,4,6-trichlorophenyl | 395 |
| pyrid-3-yloxy | 3-methylphenyl | 305 |
| pyrid-3-yloxy | 2-fluoro-5-(trifluoromethyl)phenyl | 377 |
| pyrid-3-yloxy | 3-methoxyphenyl | 321 |
| pyrid-3-yloxy | 2-bromophenyl | 370 |
| pyrid-3-yloxy | 4-bromophenyl | 370 |
| pyrid-3-yloxy | 4-fluoro-3-(trifluoromethyl)phenyl | 377 |
| pyrid-3-yloxy | 3-(trifluoromethoxy)phenyl | 375 |
| pyrid-3-yloxy | 9-fluorenon-4-yl | 393 |
| pyrid-3-yloxy | isoxazol-5-yl | 282 |
| pyrid-3-yloxy | benzofuroxan-5-yl | 349 |
| pyrid-3-yloxy | 2-chloropyrid-3-yl | 327 |
| pyrid-3-yloxy | 2-(4-methylphenoxy)pyridin-3-yl | 398 |
| pyrid-3-yloxy | pyridin-4-yl | 292 |
| pyrid-3-yloxy | anthraquinon-2-yl | 421 |
| pyrid-3-yloxy | 2-iodophenyl | 417 |
| pyrid-3-yloxy | 4-pentylphenyl | 361 |
| pyrid-3-yloxy | 2-(4-chlorophenylthio)pyridin-3-yl | 435 |
| pyrid-3-yloxy | 2,6-dimethylphenyl | 319 |
| pyrid-3-yloxy | 2,5-dimethoxyphenyl | 354 |
| pyrid-3-yloxy | 2,5-dichloropyridin-3-yl | 361 |

| | | |
|----------------|------------------------------------|-----|
| pyrid-3-yloxy | 2-chloro-6-methoxypyridin-4-yl | 357 |
| pyrid-3-yloxy | 2,3-dichloropyridin-5-yl | 361 |
| pyrid-3-yloxy | 1-naphthyl | 355 |
| pyrid-3-yloxy | 2,4-dimethoxyphenyl | 351 |
| pyrid-3-yloxy | 3,5-bis(trifluoromethyl)phenyl | 427 |
| pyrid-3-yloxy | 2-(4-chlorophenoxy)pyridin-3-yl | 419 |
| pyrid-3-yloxy | pentafluorophenyl | 381 |
| 4-bromophenoxy | 3,4-dimethoxyphenyl | 429 |
| 4-bromophenoxy | 2-(trifluoromethyl)phenyl | 437 |
| 4-bromophenoxy | 2,4-difluorophenyl | 405 |
| 4-bromophenoxy | 3-(trifluoromethyl)phenyl | 437 |
| 4-bromophenoxy | 2-naphthyl | 419 |
| 4-bromophenoxy | 2-methoxyphenyl | 399 |
| 4-bromophenoxy | 3,4,5-trimethylphenyl | 459 |
| 4-bromophenoxy | 3,4-dichlorophenyl | 438 |
| 4-bromophenoxy | 3-bromophenyl | 448 |
| 4-bromophenoxy | 3-pyridyl | 370 |
| 4-bromophenoxy | 2-ethoxynaphth-1-yl | 463 |
| 4-bromophenoxy | 2,3-dichlorophenyl | 438 |
| 4-bromophenoxy | 6-chloropyrid-3-yl | 405 |
| 4-bromophenoxy | 4-(trifluoromethoxy)phenyl | 453 |
| 4-bromophenoxy | 2-fluoro-4-(trifluoromethyl)phenyl | 455 |
| 4-bromophenoxy | 3-bromothienyl | 454 |
| 4-bromophenoxy | 2-acetoxyphenyl | 427 |

| | | |
|----------------|---|-----|
| 4-bromophenoxy | 5-methylisoxazol-3-yl | 374 |
| 4-bromophenoxy | 2-(phenylthio)pyrid-3-yl | 478 |
| 4-bromophenoxy | 2-(trifluoromethoxy)phenyl | 453 |
| 4-bromophenoxy | 1-phenyl-5-propylpyrazin-4-yl | 477 |
| 4-bromophenoxy | 2-ethoxyphenyl | 413 |
| 4-bromophenoxy | 3-chlorothien-2-yl | 410 |
| 4-bromophenoxy | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 429 |
| 4-bromophenoxy | 3,5-dichlorophenyl | 438 |
| 4-bromophenoxy | 2-(propylthio)pyridin-3-yl | 444 |
| 4-bromophenoxy | 2-(ethylthio)pyridin-3-yl | 430 |
| 4-bromophenoxy | 3-bromopyridin-5-yl | 449 |
| 4-bromophenoxy | 4-methyl-1,2,3-thiadiazol-5-yl | 391 |
| 4-bromophenoxy | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 429 |
| 4-bromophenoxy | 3-chlorobenzo[b]thiophen-2-yl | 460 |
| 4-bromophenoxy | 4-chlorophenyl | 404 |
| 4-bromophenoxy | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 450 |
| 4-bromophenoxy | benzo[b]thiophen-2-yl | 425 |
| 4-bromophenoxy | 3,4-dimethylphenyl | 397 |
| 4-bromophenoxy | 2-(phenoxy)pyridin-3-yl | 462 |
| 4-bromophenoxy | 2-(methylthio)pyridin-3-yl | 416 |
| 4-bromophenoxy | 5-methyl-3-phenylisoxazol-4-yl | 450 |

| | | |
|----------------|--|-----|
| 4-bromophenoxy | 4-chloro-1,3-dimethyl pyrazolo[3,4-b]pyridin-3- yl | 473 |
| 4-bromophenoxy | 2-chloro-6-methylpyridin- 4-yl | 419 |
| 4-bromophenoxy | 3,5-dimethylisoxazol-4-yl | 388 |
| 4-bromophenoxy | 1-naphthyl | 419 |
| 4-bromophenoxy | 2-fluorophenyl | 387 |
| 4-bromophenoxy | 4-propylphenyl | 411 |
| 4-bromophenoxy | 3-fluorophenyl | 387 |
| 4-bromophenoxy | 2,6-difluorophenyl | 405 |
| 4-bromophenoxy | 2-chlorophenyl | 414 |
| 4-bromophenoxy | 3-(chloromethyl)phenyl | 418 |
| 4-bromophenoxy | 4-(2-(2- methyl)propyl)phenyl | 425 |
| 4-bromophenoxy | 3-chlorophenyl | 404 |
| 4-bromophenoxy | 3,5-dimethoxyphenyl | 429 |
| 4-bromophenoxy | 2,6-dichlorophenyl | 438 |
| 4-bromophenoxy | 2,4-dichlorophenyl | 438 |
| 4-bromophenoxy | 4-fluorophenyl | 387 |
| 4-bromophenoxy | 4-butylphenyl | 425 |
| 4-bromophenoxy | 2-methylphenyl | 383 |
| 4-bromophenoxy | phenyl | 369 |
| 4-bromophenoxy | 4-ethylphenyl | 397 |
| 4-bromophenoxy | 2,3-difluorophenyl | 405 |
| 4-bromophenoxy | 2,6-dimethoxyphenyl | 429 |
| 4-bromophenoxy | 3,4-difluorophenyl | 405 |
| 4-bromophenoxy | 2,5-difluorophenyl | 405 |

| | | |
|----------------|------------------------------------|-----|
| 4-bromophenoxy | 4-ethoxyphenyl | 413 |
| 4-bromophenoxy | 2,4,6-trichlorophenyl | 473 |
| 4-bromophenoxy | 3-methylphenyl | 383 |
| 4-bromophenoxy | 2-fluoro-5-(trifluoromethyl)phenyl | 455 |
| 4-bromophenoxy | 3-methoxyphenyl | 399 |
| 4-bromophenoxy | 2-bromophenyl | 448 |
| 4-bromophenoxy | 4-bromophenyl | 448 |
| 4-bromophenoxy | 4-fluoro-3-(trifluoromethyl)phenyl | 455 |
| 4-bromophenoxy | 3-(trifluoromethoxy)phenyl | 453 |
| 4-bromophenoxy | 9-fluorenon-4-yl | 471 |
| 4-bromophenoxy | isoxazol-5-yl | 360 |
| 4-bromophenoxy | benzofuroxan-5-yl | 427 |
| 4-bromophenoxy | 2-chloropyrid-3-yl | 360 |
| 4-bromophenoxy | 2-(4-methylphenoxy)pyridin-3-yl | 476 |
| 4-bromophenoxy | pyridin-4-yl | 370 |
| 4-bromophenoxy | anthraquinon-2-yl | 499 |
| 4-bromophenoxy | 2-iodophenyl | 495 |
| 4-bromophenoxy | 4-pentylphenyl | 439 |
| 4-bromophenoxy | 2-(4-chlorophenylthio)pyridin-3-yl | 513 |
| 4-bromophenoxy | 2,6-dimethylphenyl | 397 |
| 4-bromophenoxy | 2,5-dimethoxyphenyl | 429 |
| 4-bromophenoxy | 2,5-dichloropyridin-3-yl | 439 |
| 4-bromophenoxy | 2-chloro-6-methoxypyridin-4-yl | 435 |

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|-----------------------------|---------------------------------|-----|
| 4-bromophenoxy | 2,3-dichloropyridin-5-yl | 439 |
| 4-bromophenoxy | 1-naphthyl | 433 |
| 4-bromophenoxy | 2,4-dimethoxyphenyl | 429 |
| 4-bromophenoxy | 3,5-bis(trifluoromethyl)phenyl | 505 |
| 4-bromophenoxy | 2-(4-chlorophenoxy)pyridin-3-yl | 497 |
| 4-bromophenoxy | pentafluorophenyl | 459 |
| 4-chloro-2-methylphenylthio | 4-biphenyl | 431 |
| 4-chloro-2-methylphenylthio | 3,4-dimethoxyphenyl | 415 |
| 4-chloro-2-methylphenylthio | 2-(trifluoromethyl)phenyl | 423 |
| 4-chloro-2-methylphenylthio | 2,4-difluorophenyl | 391 |
| 4-chloro-2-methylphenylthio | 4-cyanophenyl | 380 |
| 4-chloro-2-methylphenylthio | 3-(trifluoromethyl)phenyl | 423 |
| 4-chloro-2-methylphenylthio | 3-cyanophenyl | 380 |
| 4-chloro-2-methylphenylthio | 2-naphthyl | 405 |
| 4-chloro-2-methylphenylthio | 2-methoxyphenyl | 385 |
| 4-chloro-2-methylphenylthio | 3,4,5-trimethylphenyl | 445 |
| 4-chloro-2-methylphenylthio | 4-nitrophenyl | 400 |

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|-----------------------------|------------------------------------|-----|
| 4-chloro-2-methylphenylthio | 3,4-dichlorophenyl | 424 |
| 4-chloro-2-methylphenylthio | 5-nitrofuran-2-yl | 390 |
| 4-chloro-2-methylphenylthio | 3-bromophenyl | 434 |
| 4-chloro-2-methylphenylthio | 3-pyridyl | 356 |
| 4-chloro-2-methylphenylthio | 2-ethoxynaphth-1-yl | 449 |
| 4-chloro-2-methylphenylthio | 2,3-dichlorophenyl | 424 |
| 4-chloro-2-methylphenylthio | 3-nitrophenyl | 400 |
| 4-chloro-2-methylphenylthio | 6-chloropyrid-3-yl | 390 |
| 4-chloro-2-methylphenylthio | 4-(trifluoromethoxy)phenyl | 439 |
| 4-chloro-2-methylphenylthio | 2-fluoro-4-(trifluoromethyl)phenyl | 441 |
| 4-chloro-2-methylphenylthio | 3-bromothieryl | 440 |
| 4-chloro-2-methylphenylthio | 2-acetoxyphenyl | 413 |
| 4-chloro-2-methylphenylthio | 5-methylisoxazol-3-yl | 360 |
| 4-chloro-2-methylphenylthio | 2-(phenylthio)pyrid-3-yl | 464 |
| 4-chloro-2-methylphenylthio | 2-(trifluoromethoxy)phenyl | 439 |

| | | |
|-----------------------------|---|-----|
| 4-chloro-2-methylphenylthio | 1-phenyl-5-propylpyrazin-4-yl | 463 |
| 4-chloro-2-methylphenylthio | 2-ethoxyphenyl | 399 |
| 4-chloro-2-methylphenylthio | 3-chlorothien-2-yl | 395 |
| 4-chloro-2-methylphenylthio | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 415 |
| 4-chloro-2-methylphenylthio | 3,5-dichlorophenyl | 424 |
| 4-chloro-2-methylphenylthio | 2-(propylthio)pyridin-3-yl | 430 |
| 4-chloro-2-methylphenylthio | 2-(ethylthio)pyridin-3-yl | 416 |
| 4-chloro-2-methylphenylthio | 3-bromopyridin-5-yl | 435 |
| 4-chloro-2-methylphenylthio | 4-methyl-1,2,3-thiadiazol-5-yl | 377 |
| 4-chloro-2-methylphenylthio | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 415 |
| 4-chloro-2-methylphenylthio | 3-chlorobenzo[b]thiophen-2-yl | 445 |
| 4-chloro-2-methylphenylthio | 4-chlorophenyl | 389 |
| 4-chloro-2-methylphenylthio | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 436 |
| 4-chloro-2-methylphenylthio | benzo[b]thiophen-2-yl | 411 |
| 4-chloro-2-methylphenylthio | 3,4-dimethylphenyl | 383 |

| | | |
|-----------------------------|--|-----|
| 4-chloro-2-methylphenylthio | 2-(phenoxy)pyridin-3-yl | 448 |
| 4-chloro-2-methylphenylthio | 2-(methylthio)pyridin-3-yl | 402 |
| 4-chloro-2-methylphenylthio | 5-methyl-3-phenylisoxazol-4-yl | 436 |
| 4-chloro-2-methylphenylthio | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 458 |
| 4-chloro-2-methylphenylthio | 2-chloro-6-methylpyridin-4-yl | 404 |
| 4-chloro-2-methylphenylthio | 3,5-dimethylisoxazol-4-yl | 374 |
| 4-chloro-2-methylphenylthio | 1-naphthyl | 405 |
| 4-chloro-2-methylphenylthio | 2-fluorophenyl | 373 |
| 4-chloro-2-methylphenylthio | 4-propylphenyl | 397 |
| 4-chloro-2-methylphenylthio | 4-(trifluoromethyl)phenyl | 423 |
| 4-chloro-2-methylphenylthio | 3-fluorophenyl | 373 |
| 4-chloro-2-methylphenylthio | 2,6-difluorophenyl | 391 |
| 4-chloro-2-methylphenylthio | 2-chlorophenyl | 389 |
| 4-chloro-2-methylphenylthio | 3-(chloromethyl)phenyl | 403 |
| 4-chloro-2-methylphenylthio | 4-(2-(2-methyl)propyl)phenyl | 411 |

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|-----------------------------|-----------------------|-----|
| 4-chloro-2-methylphenylthio | 3-chlorophenyl | 389 |
| 4-chloro-2-methylphenylthio | 2-nitrophenyl | 400 |
| 4-chloro-2-methylphenylthio | 3,5-dimethoxyphenyl | 415 |
| 4-chloro-2-methylphenylthio | 2,6-dichlorophenyl | 424 |
| 4-chloro-2-methylphenylthio | 2,4-dichlorophenyl | 424 |
| 4-chloro-2-methylphenylthio | 4-fluorophenyl | 373 |
| 4-chloro-2-methylphenylthio | 4-butylphenyl | 411 |
| 4-chloro-2-methylphenylthio | 2-methylphenyl | 369 |
| 4-chloro-2-methylphenylthio | phenyl | 355 |
| 4-chloro-2-methylphenylthio | 4-ethylphenyl | 383 |
| 4-chloro-2-methylphenylthio | 2,3-difluorophenyl | 391 |
| 4-chloro-2-methylphenylthio | 2,6-dimethoxyphenyl | 415 |
| 4-chloro-2-methylphenylthio | 3,4-difluorophenyl | 391 |
| 4-chloro-2-methylphenylthio | 2,5-difluorophenyl | 391 |
| 4-chloro-2-methylphenylthio | 4-ethoxyphenyl | 399 |
| 4-chloro-2-methylphenylthio | 2,4,6-trichlorophenyl | 458 |

| | | |
|-----------------------------|------------------------------------|-----|
| 4-chloro-2-methylphenylthio | 3-methylphenyl | 369 |
| 4-chloro-2-methylphenylthio | 2-fluoro-5-(trifluoromethyl)phenyl | 441 |
| 4-chloro-2-methylphenylthio | 3-methoxyphenyl | 385 |
| 4-chloro-2-methylphenylthio | thien-2-yl | 361 |
| 4-chloro-2-methylphenylthio | 2-bromophenyl | 434 |
| 4-chloro-2-methylphenylthio | 4-bromophenyl | 434 |
| 4-chloro-2-methylphenylthio | 4-fluoro-3-(trifluoromethyl)phenyl | 441 |
| 4-chloro-2-methylphenylthio | 3-(trifluoromethoxy)phenyl | 439 |
| 4-chloro-2-methylphenylthio | 9-fluorenon-4-yl | 457 |
| 4-chloro-2-methylphenylthio | isoxazol-5-yl | 346 |
| 4-chloro-2-methylphenylthio | benzofuroxan-5-yl | 413 |
| 4-chloro-2-methylphenylthio | 2-chloropyrid-3-yl | 390 |
| 4-chloro-2-methylphenylthio | 3,5-difluorophenyl | 391 |
| 4-chloro-2-methylphenylthio | 2-(4-methylphenoxy)pyridin-3-yl | 462 |
| 4-chloro-2-methylphenylthio | pyridin-4-yl | 356 |
| 4-chloro-2-methylphenylthio | anthraquinon-2-yl | 485 |

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|-----------------------------|---------------------------|-----|
| 4-chloro-2-methylphenylthio | 2-iodophenyl | 481 |
| 4-chloro-2-methylanilino | 4-biphenyl | 414 |
| 4-chloro-2-methylanilino | 3,4-dimethoxyphenyl | 398 |
| 4-chloro-2-methylanilino | 2-(trifluoromethyl)phenyl | 406 |
| 4-chloro-2-methylanilino | 2,4-difluorophenyl | 374 |
| 4-chloro-2-methylanilino | 4-cyanophenyl | 363 |
| 4-chloro-2-methylanilino | 3-(trifluoromethyl)phenyl | 406 |
| 4-chloro-2-methylanilino | 3-cyanophenyl | 363 |
| 4-chloro-2-methylanilino | 2-naphthyl | 388 |
| 4-chloro-2-methylanilino | 2-methoxyphenyl | 368 |
| 4-chloro-2-methylanilino | 3,4,5-trimethylphenyl | 428 |
| 4-chloro-2-methylanilino | 4-nitrophenyl | 383 |
| 4-chloro-2-methylanilino | 3,4-dichlorophenyl | 407 |
| 4-chloro-2-methylanilino | 5-nitrofuran-2-yl | 373 |
| 4-chloro-2-methylanilino | 3-bromophenyl | 417 |
| 4-chloro-2-methylanilino | 3-pyridyl | 339 |

| | | |
|--------------------------|---|-----|
| 4-chloro-2-methylanilino | 2-ethoxynaphth-1-yl | 432 |
| 4-chloro-2-methylanilino | 2,3-dichlorophenyl | 407 |
| 4-chloro-2-methylanilino | 3-nitrophenyl | 383 |
| 4-chloro-2-methylanilino | 6-chloropyrid-3-yl | 373 |
| 4-chloro-2-methylanilino | 4-(trifluoromethoxy)phenyl | 422 |
| 4-chloro-2-methylanilino | 2-fluoro-4-(trifluoromethyl)phenyl | 424 |
| 4-chloro-2-methylanilino | 3-bromothienyl | 423 |
| 4-chloro-2-methylanilino | 2-acetoxyphenyl | 396 |
| 4-chloro-2-methylanilino | 5-methylisoxazol-3-yl | 343 |
| 4-chloro-2-methylanilino | 2-(phenylthio)pyrid-3-yl | 447 |
| 4-chloro-2-methylanilino | 2-(trifluoromethoxy)phenyl | 422 |
| 4-chloro-2-methylanilino | 1-phenyl-5-propylpyrazin-4-yl | 446 |
| 4-chloro-2-methylanilino | 2-ethoxyphenyl | 382 |
| 4-chloro-2-methylanilino | 3-chlorothien-2-yl | 378 |
| 4-chloro-2-methylanilino | 1-(2-(2-methyl)propyl)-3-methylpyrazol-5-yl | 398 |
| 4-chloro-2-methylanilino | 3,5-dichlorophenyl | 407 |

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|--------------------------|--|-----|
| 4-chloro-2-methylanilino | 2-(propylthio)pyridin-3-yl | 413 |
| 4-chloro-2-methylanilino | 2-(ethylthio)pyridin-3-yl | 399 |
| 4-chloro-2-methylanilino | 3-bromopyridin-5-yl | 418 |
| 4-chloro-2-methylanilino | 4-methyl-1,2,3-thiadiazol-5-yl | 360 |
| 4-chloro-2-methylanilino | 1-methyl-3-(2-(2-methyl)propyl)pyrazol-5-yl | 398 |
| 4-chloro-2-methylanilino | 3-chlorobenzo[b]thiophen-2-yl | 428 |
| 4-chloro-2-methylanilino | 4-chlorophenyl | 372 |
| 4-chloro-2-methylanilino | 4-methyl-2-phenyl-1,2,3-triazol-5-yl | 419 |
| 4-chloro-2-methylanilino | benzo[b]thiophen-2-yl | 394 |
| 4-chloro-2-methylanilino | 3,4-dimethylphenyl | 366 |
| 4-chloro-2-methylanilino | 2-(phenoxy)pyridin-3-yl | 431 |
| 4-chloro-2-methylanilino | 2-(methylthio)pyridin-3-yl | 385 |
| 4-chloro-2-methylanilino | 5-methyl-3-phenylisoxazol-4-yl | 419 |
| 4-chloro-2-methylanilino | 4-chloro-1,3-dimethylpyrazolo[3,4-b]pyridin-3-yl | 441 |
| 4-chloro-2-methylanilino | 2-chloro-6-methylpyridin-4-yl | 387 |

| | | |
|--------------------------|------------------------------|-----|
| 4-chloro-2-methylanilino | 3,5-dimethylisoxazol-4-yl | 357 |
| 4-chloro-2-methylanilino | 1-naphthyl | 388 |
| 4-chloro-2-methylanilino | 2-fluorophenyl | 356 |
| 4-chloro-2-methylanilino | 4-propylphenyl | 380 |
| 4-chloro-2-methylanilino | 4-(trifluoromethyl)phenyl | 406 |
| 4-chloro-2-methylanilino | 3-fluorophenyl | 356 |
| 4-chloro-2-methylanilino | 2,6-difluorophenyl | 374 |
| 4-chloro-2-methylanilino | 2-chlorophenyl | 372 |
| 4-chloro-2-methylanilino | 3-(chloromethyl)phenyl | 386 |
| 4-chloro-2-methylanilino | 4-(2-(2-methyl)propyl)phenyl | 394 |
| 4-chloro-2-methylanilino | 3-chlorophenyl | 372 |
| 4-chloro-2-methylanilino | 2-nitrophenyl | 383 |
| 4-chloro-2-methylanilino | 3,5-dimethoxyphenyl | 398 |
| 4-chloro-2-methylanilino | 2,6-dichlorophenyl | 407 |
| 4-chloro-2-methylanilino | 2,4-dichlorophenyl | 407 |

| | | |
|--------------------------|------------------------------------|-----|
| 4-chloro-2-methylanilino | 4-fluorophenyl | 356 |
| 4-chloro-2-methylanilino | 4-butylphenyl | 394 |
| 4-chloro-2-methylanilino | 2-methylphenyl | 352 |
| 4-chloro-2-methylanilino | phenyl | 338 |
| 4-chloro-2-methylanilino | 4-ethylphenyl | 366 |
| 4-chloro-2-methylanilino | 2,3-difluorophenyl | 374 |
| 4-chloro-2-methylanilino | 2,6-dimethoxyphenyl | 398 |
| 4-chloro-2-methylanilino | 3,4-difluorophenyl | 374 |
| 4-chloro-2-methylanilino | 2,5-difluorophenyl | 374 |
| 4-chloro-2-methylanilino | 4-ethoxyphenyl | 382 |
| 4-chloro-2-methylanilino | 2,4,6-trichlorophenyl | 441 |
| 4-chloro-2-methylanilino | 3-methylphenyl | 352 |
| 4-chloro-2-methylanilino | 2-fluoro-5-(trifluoromethyl)phenyl | 424 |
| 4-chloro-2-methylanilino | 3-methoxyphenyl | 368 |
| 4-chloro-2-methylanilino | thien-2-yl | 344 |

| | | |
|--------------------------|------------------------------------|-----|
| 4-chloro-2-methylanilino | 2-bromophenyl | 417 |
| 4-chloro-2-methylanilino | 4-bromophenyl | 417 |
| 4-chloro-2-methylanilino | 4-fluoro-3-(trifluoromethyl)phenyl | 424 |
| 4-chloro-2-methylanilino | 3-(trifluoromethoxy)phenyl | 422 |
| 4-chloro-2-methylanilino | 9-fluorenon-4-yl | 440 |
| 4-chloro-2-methylanilino | isoxazol-5-yl | 329 |
| 4-chloro-2-methylanilino | benzofuroxan-5-yl | 396 |
| 4-chloro-2-methylanilino | 2-chloropyrid-3-yl | 373 |
| 4-chloro-2-methylanilino | 3,5-difluorophenyl | 374 |
| 4-chloro-2-methylanilino | 2-(4-methylphenoxy)pyridin-3-yl | 445 |
| 4-chloro-2-methylanilino | pyridin-4-yl | 339 |
| 4-chloro-2-methylanilino | anthraquinon-2-yl | 468 |
| 4-chloro-2-methylanilino | 2-iodophenyl | 464 |

The compounds listed in Table 7 can be prepared from substituted 5-aminopyridine compounds and the appropriate acid chloride according to the general procedure above.

Table 7

| <u>R¹X</u> | <u>R³</u> |
|--------------------------|--|
| 4-chloro-2-methylphenoxy | 3,4-difluorophenyl |
| 4-chloro-2-methylphenoxy | 4-pentylphenyl |
| 4-chloro-2-methylphenoxy | 2-(4-chlorophenylthio) pyridin-3-yl |
| 4-chloro-2-methylphenoxy | 2,6-dimethylphenyl |
| 4-chloro-2-methylphenoxy | 2,5-dimethoxyphenyl |
| 4-chloro-2-methylphenoxy | 2,5-dichloropyridin-3-yl |
| 4-chloro-2-methylphenoxy | 2-chloro-6-methoxypyridin- 4-yl |
| 4-chloro-2-methylphenoxy | 2,3-dichloropyridin-5-yl |
| 4-chloro-2-methylphenoxy | 1-naphthyl |
| 4-chloro-2-methylphenoxy | 2,4-dimethoxyphenyl |
| 4-chloro-2-methylphenoxy | 3,5- bis(trifluoromethyl)phenyl |
| 4-chloro-2-methylphenoxy | 2-(4- chlorophenoxy)pyridin-3-yl |
| 4-chloro-2-methylphenoxy | pentafluorophenyl |
| 1-naphthoxy | 4-pentylphenyl |
| 1-naphthoxy | 2-(4-chlorophenylthio) pyridin-3-yl |
| 1-naphthoxy | 2,6-dimethylphenyl |
| 1-naphthoxy | 2,5-dimethoxyphenyl |
| 1-naphthoxy | 2,5-dichloropyridin-3-yl |
| 1-naphthoxy | 2-chloro-6-methoxypyridin- 4-yl |
| 1-naphthoxy | 2,3-dichloropyridin-5-yl |
| 1-naphthoxy | 1-naphthyl |

| | |
|--------------------------|--|
| 1-naphthoxy | 2,4-dimethoxyphenyl |
| 1-naphthoxy | 3,5-bis(trifluoromethyl) phenyl |
| 1-naphthoxy | 2-(4- chlorophenoxy)pyridin-3-yl |
| 1-naphthoxy | pentafluorophenyl |
| 2-(2-propyl)phenoxy | 4-pentylphenyl |
| 2-(2-propyl)phenoxy | 2-(4-chlorophenylthio) pyridin-3-yl |
| 2-(2-propyl)phenoxy | 2,6-dimethylphenyl |
| 2-(2-propyl)phenoxy | 2,5-dimethoxyphenyl |
| 2-(2-propyl)phenoxy | 2,5-dichloropyridin-3-yl |
| 2-(2-propyl)phenoxy | 2-chloro-6-methoxypyridin- 4-yl |
| 2-(2-propyl)phenoxy | 2,3-dichloropyridin-5-yl |
| 2-(2-propyl)phenoxy | 1-naphthyl |
| 2-(2-propyl)phenoxy | 2,4-dimethoxyphenyl |
| 2-(2-propyl)phenoxy | 3,5-bis(trifluoromethyl) phenyl |
| 2-(2-propyl)phenoxy | 2-(4- chlorophenoxy)pyridin-3-yl |
| 2-(2-propyl)phenoxy | pentafluorophenyl |
| 3-fluoro-5-methylphenoxy | 4-pentylphenyl |
| 3-fluoro-5-methylphenoxy | 2-(4-chlorophenylthio) pyridin-3-yl |
| 3-fluoro-5-methylphenoxy | 2,6-dimethylphenyl |
| 3-fluoro-5-methylphenoxy | 2,5-dimethoxyphenyl |
| 3-fluoro-5-methylphenoxy | 2,5-dichloropyridin-3-yl |
| 3-fluoro-5-methylphenoxy | 2-chloro-6-methoxypyridin- 4-yl |

| | |
|--------------------------|--|
| 3-fluoro-5-methylphenoxy | 2,3-dichloropyridin-5-yl |
| 3-fluoro-5-methylphenoxy | 1-naphthyl |
| 3-fluoro-5-methylphenoxy | 2,4-dimethoxyphenyl |
| 3-fluoro-5-methylphenoxy | 3,5-bis(trifluoromethyl) phenyl |
| 3-fluoro-5-methylphenoxy | 2-(4- chlorophenoxy)pyridin-3-yl |
| 3-fluoro-5-methylphenoxy | pentafluorophenyl |
| 2-methylpyrid-3-yloxy | 4-pentylphenyl |
| 2-methylpyrid-3-yloxy | 2-(4-chlorophenylthio) pyridin-3-yl |
| 2-methylpyrid-3-yloxy | 2,6-dimethylphenyl |
| 2-methylpyrid-3-yloxy | 2,5-dimethoxyphenyl |
| 2-methylpyrid-3-yloxy | 2,5-dichloropyridin-3-yl |
| 2-methylpyrid-3-yloxy | 2-chloro-6-methoxypyridin- 4-yl |
| 2-methylpyrid-3-yloxy | 2,3-dichloropyridin-5-yl |
| 2-methylpyrid-3-yloxy | 1-naphthyl |
| 2-methylpyrid-3-yloxy | 2,4-dimethoxyphenyl |
| 2-methylpyrid-3-yloxy | 3,5-bis(trifluoromethyl) phenyl |
| 2-methylpyrid-3-yloxy | 2-(4- chlorophenoxy)pyridin-3-yl |
| 2-methylpyrid-3-yloxy | pentafluorophenyl |
| 4-methoxyphenoxy | 4-biphenyl |
| 4-methoxyphenoxy | 4-cyanophenyl |
| 4-methoxyphenoxy | 3-cyanophenyl |
| 4-methoxyphenoxy | 4-nitrophenyl |
| 4-methoxyphenoxy | 5-nitrofuran-2-yl |

| | |
|----------------------|---------------------------|
| 4-methoxyphenoxy | 3-nitrophenyl |
| 4-methoxyphenoxy | 4-(trifluoromethyl)phenyl |
| 4-methoxyphenoxy | 2-nitrophenyl |
| 4-methoxyphenoxy | thien-2-yl |
| 2-(2-propoxy)phenoxy | 4-biphenyl |
| 2-(2-propoxy)phenoxy | 4-cyanophenyl |
| 2-(2-propoxy)phenoxy | 3-cyanophenyl |
| 2-(2-propoxy)phenoxy | 4-nitrophenyl |
| 2-(2-propoxy)phenoxy | 5-nitrofuran-2-yl |
| 2-(2-propoxy)phenoxy | 3-nitrophenyl |
| 2-(2-propoxy)phenoxy | 2-nitrophenyl |
| 2-(2-propoxy)phenoxy | thien-2-yl |
| 2-(2-propoxy)phenoxy | 3,5-difluorophenyl |
| 4-fluorophenoxy | 4-biphenyl |
| 4-fluorophenoxy | 4-cyanophenyl |
| 4-fluorophenoxy | 3-cyanophenyl |
| 4-fluorophenoxy | 4-nitrophenyl |
| 4-fluorophenoxy | 5-nitrofuran-2-yl |
| 4-fluorophenoxy | 3-nitrophenyl |
| 4-fluorophenoxy | 2-nitrophenyl |
| 4-fluorophenoxy | 4-(trifluoromethyl)phenyl |
| 4-fluorophenoxy | thien-2-yl |
| 4-fluorophenoxy | 3,5-difluorophenyl |
| 4-chlorophenoxy | 4-biphenyl |
| 4-chlorophenoxy | 4-cyanophenyl |
| 4-chlorophenoxy | 3-cyanophenyl |
| 4-chlorophenoxy | 4-nitrophenyl |
| 4-chlorophenoxy | 5-nitrofuran-2-yl |

| | |
|------------------------------|---------------------------|
| 4-chlorophenoxy | 3-nitrophenyl |
| 4-chlorophenoxy | 2-nitrophenyl |
| 4-chlorophenoxy | 4-(trifluoromethyl)phenyl |
| 4-chlorophenoxy | thien-2-yl |
| 4-chlorophenoxy | 3,5-difluorophenyl |
| 2,4-difluorophenoxy | 4-biphenyl |
| 2,4-difluorophenoxy | 4-cyanophenyl |
| 2,4-difluorophenoxy | 3-cyanophenyl |
| 2,4-difluorophenoxy | 4-nitrophenyl |
| 2,4-difluorophenoxy | 5-nitrofuran-2-yl |
| 2,4-difluorophenoxy | 3-nitrophenyl |
| 2,4-difluorophenoxy | 4-(trifluoromethyl)phenyl |
| 2,4-difluorophenoxy | 2-nitrophenyl |
| 4-chloro-2,5-dimethylphenoxy | 4-biphenyl |
| 4-chloro-2,5-dimethylphenoxy | 4-cyanophenyl |
| 4-chloro-2,5-dimethylphenoxy | 3-cyanophenyl |
| 4-chloro-2,5-dimethylphenoxy | 4-nitrophenyl |
| 4-chloro-2,5-dimethylphenoxy | 5-nitrofuran-2-yl |
| 4-chloro-2,5-dimethylphenoxy | 3-nitrophenyl |
| 4-chloro-2,5-dimethylphenoxy | 4-(trifluoromethyl)phenyl |
| 4-chloro-2,5-dimethylphenoxy | 2-nitrophenyl |

| | |
|-------------------------------|---------------------------|
| 4-chloro-2,5-dimethylphenoxy | thien-2-yl |
| 4-chloro-2,5-dimethylphenoxy | 3,5-difluorophenyl |
| 4-methoxyphenoxy | 3,5-difluorophenyl |
| 2-(2-propoxy)phenoxy | 4-(trifluoromethyl)phenyl |
| 2,4-difluorophenoxy | thien-2-yl |
| 2,4-difluorophenoxy | 3,5-difluorophenyl |
| 4-thiomethylphenoxy | 4-biphenyl |
| 4-thiomethylphenoxy | 4-cyanophenyl |
| 4-thiomethylphenoxy | 3-cyanophenyl |
| 4-thiomethylphenoxy | 4-nitrophenyl |
| 4-thiomethylphenoxy | 5-nitrofuran-2-yl |
| 4-thiomethylphenoxy | 3-nitrophenyl |
| 4-thiomethylphenoxy | 4-(trifluoromethyl)phenyl |
| 4-thiomethylphenoxy | 2-nitrophenyl |
| 4-thiomethylphenoxy | thien-2-yl |
| 4-thiomethylphenoxy | 3,5-difluorophenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 4-biphenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 4-cyanophenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 3-cyanophenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 4-nitrophenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 5-nitrofuran-2-yl |

| | |
|-------------------------------|---------------------------|
| 4-(2-(2-methyl)propyl)phenoxy | 3-nitrophenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 4-(trifluoromethyl)phenyl |
| 4-(2-(2-methyl)propyl)phenoxy | 2-nitrophenyl |
| 4-(2-(2-methyl)propyl)phenoxy | thien-2-yl |
| 4-(2-(2-methyl)propyl)phenoxy | 3,5-difluorophenyl |
| 2,3-dimethylphenoxy | 4-biphenyl |
| 2,3-dimethylphenoxy | 4-cyanophenyl |
| 2,3-dimethylphenoxy | 3-cyanophenyl |
| 2,3-dimethylphenoxy | 4-nitrophenyl |
| 2,3-dimethylphenoxy | 5-nitrofuran-2-yl |
| 2,3-dimethylphenoxy | 3-nitrophenyl |
| 2,3-dimethylphenoxy | 4-(trifluoromethyl)phenyl |
| 2,3-dimethylphenoxy | 2-nitrophenyl |
| 2,3-dimethylphenoxy | thien-2-yl |
| 2,3-dimethylphenoxy | 3,5-difluorophenyl |
| 3,5-(bis-2-propyl)phenoxy | 4-biphenyl |
| 3,5-(bis-2-propyl)phenoxy | 4-cyanophenyl |
| 3,5-(bis-2-propyl)phenoxy | 3-cyanophenyl |
| 3,5-(bis-2-propyl)phenoxy | 4-nitrophenyl |
| 3,5-(bis-2-propyl)phenoxy | 5-nitrofuran-2-yl |
| 3,5-(bis-2-propyl)phenoxy | 3-nitrophenyl |
| 3,5-(bis-2-propyl)phenoxy | 4-(trifluoromethyl)phenyl |
| 3,5-(bis-2-propyl)phenoxy | 2-nitrophenyl |

| | |
|---------------------------|---------------------------|
| 3,5-(bis-2-propyl)phenoxy | thien-2-yl |
| 3,5-(bis-2-propyl)phenoxy | 3,5-difluorophenyl |
| 3-trifluoromethyl phenoxy | 4-biphenyl |
| 3-trifluoromethyl phenoxy | 4-cyanophenyl |
| 3-trifluoromethyl phenoxy | 3-cyanophenyl |
| 3-trifluoromethyl phenoxy | 4-nitrophenyl |
| 3-trifluoromethyl phenoxy | 5-nitrofuran-2-yl |
| 3-trifluoromethyl phenoxy | 3-nitrophenyl |
| 3-trifluoromethyl phenoxy | 4-(trifluoromethyl)phenyl |
| 3-trifluoromethyl phenoxy | 2-nitrophenyl |
| 3-trifluoromethyl phenoxy | thien-2-yl |
| 3-trifluoromethyl phenoxy | 3,5-difluorophenyl |
| 2,6-dichlorophenoxy | 4-biphenyl |
| 2,6-dichlorophenoxy | 4-cyanophenyl |
| 2,6-dichlorophenoxy | 3-cyanophenyl |
| 2,6-dichlorophenoxy | 4-nitrophenyl |
| 2,6-dichlorophenoxy | 5-nitrofuran-2-yl |
| 2,6-dichlorophenoxy | 3-nitrophenyl |
| 2,6-dichlorophenoxy | 4-(trifluoromethyl)phenyl |
| 2,6-dichlorophenoxy | 2-nitrophenyl |
| 2,6-dichlorophenoxy | thien-2-yl |
| 2,6-dichlorophenoxy | 3,5-difluorophenyl |
| 2,4-dichlorophenoxy | 4-biphenyl |
| 2,4-dichlorophenoxy | 4-cyanophenyl |
| 2,4-dichlorophenoxy | 3-cyanophenyl |
| 2,4-dichlorophenoxy | 4-nitrophenyl |
| 2,4-dichlorophenoxy | 5-nitrofuran-2-yl |
| 2,4-dichlorophenoxy | 3-nitrophenyl |

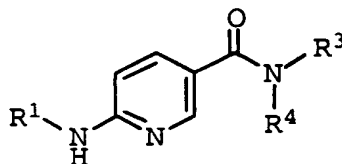
| | |
|------------------------------|---------------------------|
| 2,4-dichlorophenoxy | 4-(trifluoromethyl)phenyl |
| 2,4-dichlorophenoxy | 2-nitrophenyl |
| 2,4-dichlorophenoxy | thien-2-yl |
| 2,4-dichlorophenoxy | 3,5-difluorophenyl |
| 4-chloro-3-methylphenoxy | 4-biphenyl |
| 4-chloro-3-methylphenoxy | 4-cyanophenyl |
| 4-chloro-3-methylphenoxy | 3-cyanophenyl |
| 4-chloro-3-methylphenoxy | 4-nitrophenyl |
| 4-chloro-3-methylphenoxy | 5-nitrofuran-2-yl |
| 4-chloro-3-methylphenoxy | 3-nitrophenyl |
| 4-chloro-3-methylphenoxy | 2-nitrophenyl |
| 4-chloro-3-methylphenoxy | thien-2-yl |
| 4-chloro-3-methylphenoxy | 3,5-difluorophenyl |
| 4-chloro-2-cyclohexylphenoxy | 4-biphenyl |
| 4-chloro-2-cyclohexylphenoxy | 4-cyanophenyl |
| 4-chloro-2-cyclohexylphenoxy | 3-cyanophenyl |
| 4-chloro-2-cyclohexylphenoxy | 4-nitrophenyl |
| 4-chloro-2-cyclohexylphenoxy | 5-nitrofuran-2-yl |
| 4-chloro-2-cyclohexylphenoxy | 3-nitrophenyl |
| 4-chloro-2-cyclohexylphenoxy | 4-(trifluoromethyl)phenyl |
| 4-chloro-2-cyclohexylphenoxy | 2-nitrophenyl |

| | |
|------------------------------|---------------------------|
| 4-chloro-2-cyclohexylphenoxy | thien-2-yl |
| 4-chloro-2-cyclohexylphenoxy | 3,5-difluorophenyl |
| 4-chloro-3,5-dimethylphenoxy | 4-biphenyl |
| 4-chloro-3,5-dimethylphenoxy | 4-cyanophenyl |
| 4-chloro-3,5-dimethylphenoxy | 3-cyanophenyl |
| 4-chloro-3,5-dimethylphenoxy | 4-nitrophenyl |
| 4-chloro-3,5-dimethylphenoxy | 5-nitrofuran-2-yl |
| 4-chloro-3,5-dimethylphenoxy | 3-nitrophenyl |
| 4-chloro-3,5-dimethylphenoxy | 4-(trifluoromethyl)phenyl |
| 4-chloro-3,5-dimethylphenoxy | 2-nitrophenyl |
| 4-chloro-3,5-dimethylphenoxy | thien-2-yl |
| 4-chloro-3,5-dimethylphenoxy | 3,5-difluorophenyl |
| pyrid-3-yloxy | 4-biphenyl |
| pyrid-3-yloxy | 4-cyanophenyl |
| pyrid-3-yloxy | 3-cyanophenyl |
| pyrid-3-yloxy | 4-nitrophenyl |
| pyrid-3-yloxy | 5-nitrofuran-2-yl |
| pyrid-3-yloxy | 3-nitrophenyl |

| | |
|-----------------------------|------------------------------------|
| pyrid-3-yloxy | 4-(trifluoromethyl)phenyl |
| pyrid-3-yloxy | 2-nitrophenyl |
| pyrid-3-yloxy | thien-2-yl |
| pyrid-3-yloxy | 3,5-difluorophenyl |
| 4-bromophenoxy | 4-biphenyl |
| 4-bromophenoxy | 4-cyanophenyl |
| 4-bromophenoxy | 3-cyanophenyl |
| 4-bromophenoxy | 4-nitrophenyl |
| 4-bromophenoxy | 5-nitrofuran-2-yl |
| 4-bromophenoxy | 3-nitrophenyl |
| 4-bromophenoxy | 4-(trifluoromethyl)phenyl |
| 4-bromophenoxy | 2-nitrophenyl |
| 4-bromophenoxy | thien-2-yl |
| 4-bromophenoxy | 3,5-difluorophenyl |
| 4-chloro-2-methylphenylthio | 4-pentylphenyl |
| 4-chloro-2-methylphenylthio | 2-(4-chlorophenylthio)pyridin-3-yl |
| 4-chloro-2-methylphenylthio | 2,6-dimethylphenyl |
| 4-chloro-2-methylphenylthio | 2,5-dimethoxyphenyl |
| 4-chloro-2-methylphenylthio | 2,5-dichloropyridin-3-yl |
| 4-chloro-2-methylphenylthio | 2-chloro-6-methoxypyridin-4-yl |
| 4-chloro-2-methylphenylthio | 2,3-dichloropyridin-5-yl |
| 4-chloro-2-methylphenylthio | 1-naphthyl |

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| | |
|-----------------------------|------------------------------------|
| 4-chloro-2-methylphenylthio | 2,4-dimethoxyphenyl |
| 4-chloro-2-methylphenylthio | 3,5-bis(trifluoromethyl)phenyl |
| 4-chloro-2-methylphenylthio | 2-(4-chlorophenoxy)pyridin-3-yl |
| 4-chloro-2-methylphenylthio | pentafluorophenyl |
| 4-chloro-2-methylanilino | 4-pentylphenyl |
| 4-chloro-2-methylanilino | 2-(4-chlorophenylthio)pyridin-3-yl |
| 4-chloro-2-methylanilino | 2,6-dimethylphenyl |
| 4-chloro-2-methylanilino | 2,5-dimethoxyphenyl |
| 4-chloro-2-methylanilino | 2,5-dichloropyridin-3-yl |
| 4-chloro-2-methylanilino | 2-chloro-6-methoxypyridin-4-yl |
| 4-chloro-2-methylanilino | 2,3-dichloropyridin-5-yl |
| 4-chloro-2-methylanilino | 1-naphthyl |
| 4-chloro-2-methylanilino | 2,4-dimethoxyphenyl |
| 4-chloro-2-methylanilino | 3,5-bis(trifluoromethyl)phenyl |
| 4-chloro-2-methylanilino | 2-(4-chlorophenoxy)pyridin-3-yl |
| 4-chloro-2-methylanilino | pentafluorophenyl |

Example 19

General procedure for the synthesis of 6-(substituted-amino)-N-substituted nicotinamides

Step A. General procedure for the preparation of 6-chloro-N-substituted nicotinamide:

To a suspension of 6-chloronicotinoyl chloride (1.76 g, 10.0 mmol) in dry dichloromethane (10 mL) was added the amine (R^1R^2NH) (10.0 mmol) followed by the dropwise addition of triethylamine (1.7 mL, 12.2 mmol). After stirring for 40 min. at room temperature, the mixture was diluted with dichloromethane, washed with aqueous 1 M hydrochloric acid, saturated aqueous sodium hydrogencarbonate and water, dried over sodium sulfate and concentrated to dryness under reduce pressure to afford the desired nicotinamide.

The following compounds were prepared according to this procedure using the appropriate substituted amine:

- 6-Chloro-N-o-tolylnicotinamide: MS (m/z): 247/249 ($M+H$)⁺; $C_{13}H_{11}Cl_1N_2O_1$ requires 246.5.
- 6-Chloro-N-(2-fluorophenyl)nicotinamide: MS (m/z): 251/253 ($M+H$)⁺; $C_{12}H_8Cl_1F_1N_2O_1$ requires 250.7.
- 6-Chloro-N-(2,6-dimethylphenyl)nicotinamide: MS (m/z): 261/263 ($M+H$)⁺; $C_{14}H_{13}Cl_1N_2O_1$ requires 260.7.
- 6-Chloro-N-(2-phenoxyphenyl)nicotinamide: MS (m/z): 325/327 ($M+H$)⁺; $C_{18}H_{13}Cl_1N_2O_1$ requires 324.8.
- 6-Chloro-N-phenylnicotinamide: MS (m/z): 233/235 ($M+H$)⁺; $C_{12}H_9Cl_1N_2O_1$ requires 232.7.
- 6-Chloro-N-(2,4-difluorophenyl)nicotinamide: MS (m/z): 269/271 ($M+H$)⁺; $C_{12}H_7Cl_1F_2N_2O_1$ requires 268.6.
- 6-Chloro-N-(2,6-diisopropylphenyl)nicotinamide: MS (m/z): 317/319 ($M+H$)⁺; $C_{18}H_{21}Cl_1N_2O_1$ requires 316.8.
- 6-Chloro-N-(4-chlorophenyl)-N-methylnicotinamide: MS (m/z): 281/283 ($M+H$)⁺; $C_{13}H_{10}Cl_2N_2O_1$ requires 281.1.
- 6-Chloro-N-(2,4-dimethoxyphenyl)nicotinamide: MS (m/z): 293/295 ($M+H$)⁺; $C_{14}H_{13}Cl_1N_2O_3$ requires 292.7.
- 6-Chloro-N-(3-methoxyphenyl)nicotinamide: MS (m/z): 263/265 ($M+H$)⁺; $C_{13}H_{11}Cl_1N_2O_2$ requires 262.7.
- 6-Chloro-N-(4-methoxyphenyl)nicotinamide: MS (m/z): 263/265 ($M+H$)⁺; $C_{13}H_{11}Cl_1N_2O_2$ requires 262.7.

6-Chloro-N-(2-methoxyphenyl)nicotinamide: MS (m/z):
263/265 ($M+H$)⁺; $C_{13}H_{11}Cl_1N_2O_2$ requires 262.7.

6-Chloro-N-methyl-N-phenylnicotinamide: MS (m/z):
247/249 ($M+H$)⁺; $C_{13}H_{11}Cl_1N_2O_1$ requires 246.7.

- 5 N-Benzyl-6-chloronicotinamide: MS (m/z): 247/249 ($M+H$)⁺;
 $C_{13}H_{11}Cl_1N_2O_1$ requires 246.7.

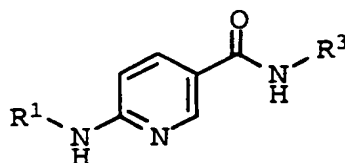
Step B. General procedure for the preparation of 6-
(substituted-amino)-N-substituted nicotinamides

- 10 A mixture of the 6-chloro-N-substituted nicotinamide
(12.5 mmol) and amine (R^1NH_2 or R^1NHCH_3) (20 mmol) in
ethylene glycol (50 mL) or pyridine (alkylamines) (50
mL) was heated to 140°C for 20 hours. After cooling to
room temperature, the mixture was diluted with
15 dichloromethane/methanol (9:1, 250 mL) and filtered
through a plug of silica gel, washing with additional
dichloromethane/methanol (9:1, 250 mL). Concentration
under reduced pressure afforded the desired 6-
(substituted-amino)-N-substituted nicotinamide.

- 20 The compounds listed in Tables 8-11 were prepared
from 6-chloro-N-substituted nicotinamides compounds and
the appropriate amine according to the general procedure
above.

25

Table 8



| R^3 | R^1 | <u>MS</u> (m/z) |
|---------|-------------------------|------------------------|
| o-tolyl | phenyl | 303 |
| o-tolyl | o-tolyl | 317 |
| o-tolyl | 4-chloro-2-methylphenyl | 352 |
| o-tolyl | 2-fluorophenyl | 321 |

| | | |
|----------------|----------------------------|-----|
| o-tolyl | 3-fluorophenyl | 321 |
| o-tolyl | 4-fluorophenyl | 321 |
| o-tolyl | 2,4-difluorophenyl | 339 |
| o-tolyl | 2-methoxyphenyl | 333 |
| o-tolyl | 3-methoxyphenyl | 333 |
| o-tolyl | 4-methoxyphenyl | 333 |
| o-tolyl | 2,4-dimethoxyphenyl | 363 |
| o-tolyl | 2-phenoxyphenyl | 395 |
| o-tolyl | 3-phenoxyphenyl | 395 |
| o-tolyl | 4-phenoxyphenyl | 395 |
| o-tolyl | 4-biphenyl | 379 |
| o-tolyl | 4-benzylphenyl | 393 |
| o-tolyl | 4-(trifluoromethoxy)phenyl | 387 |
| o-tolyl | cyclohexyl | 309 |
| o-tolyl | 2-methylcyclohexyl | 323 |
| o-tolyl | cycloheptyl | 323 |
| o-tolyl | indan-1-yl | 343 |
| o-tolyl | 2-dicyclohexyl | 492 |
| 2-fluorophenyl | phenyl | 307 |
| 2-fluorophenyl | o-tolyl | 321 |
| 2-fluorophenyl | 4-chloro-2-methylphenyl | 356 |
| 2-fluorophenyl | 2-fluorophenyl | 325 |
| 2-fluorophenyl | 3-fluorophenyl | 325 |
| 2-fluorophenyl | 4-fluorophenyl | 325 |
| 2-fluorophenyl | 2,4-difluorophenyl | 343 |
| 2-fluorophenyl | 2-methoxyphenyl | 337 |
| 2-fluorophenyl | 3-methoxyphenyl | 337 |
| 2-fluorophenyl | 4-methoxyphenyl | 337 |

| | | |
|--------------------|----------------------------|-----|
| 2-fluorophenyl | 2,4-dimethoxyphenyl | 367 |
| 2-fluorophenyl | 2-phenoxyphenyl | 399 |
| 2-fluorophenyl | 3-phenoxyphenyl | 399 |
| 2-fluorophenyl | 4-phenoxyphenyl | 399 |
| 2-fluorophenyl | 4-biphenyl | 383 |
| 2-fluorophenyl | 4-benzylphenyl | 397 |
| 2-fluorophenyl | 4-(trifluoromethoxy)phenyl | 391 |
| 2-fluorophenyl | cyclohexyl | 313 |
| 2-fluorophenyl | 2-methylcyclohexyl | 327 |
| 2-fluorophenyl | cycloheptyl | 327 |
| 2-fluorophenyl | indan-1-yl | 347 |
| 2-fluorophenyl | 2-dicyclohexyl | 395 |
| 2,6-dimethylphenyl | phenyl | 317 |
| 2,6-dimethylphenyl | o-tolyl | 331 |
| 2,6-dimethylphenyl | 4-chloro-2-methylphenyl | 366 |
| 2,6-dimethylphenyl | 2-fluorophenyl | 335 |
| 2,6-dimethylphenyl | 3-fluorophenyl | 335 |
| 2,6-dimethylphenyl | 4-fluorophenyl | 335 |
| 2,6-dimethylphenyl | 2,4-difluorophenyl | 353 |
| 2,6-dimethylphenyl | 2-methoxyphenyl | 347 |
| 2,6-dimethylphenyl | 3-methoxyphenyl | 347 |
| 2,6-dimethylphenyl | 4-methoxyphenyl | 347 |
| 2,6-dimethylphenyl | 2,4-dimethoxyphenyl | 377 |
| 2,6-dimethylphenyl | 2-phenoxyphenyl | 409 |
| 2,6-dimethylphenyl | 3-phenoxyphenyl | 409 |
| 2,6-dimethylphenyl | 4-phenoxyphenyl | 409 |
| 2,6-dimethylphenyl | 4-biphenyl | 393 |
| 2,6-dimethylphenyl | 4-benzylphenyl | 407 |

| | | |
|--------------------|----------------------------|-----|
| 2,6-dimethylphenyl | 4-(trifluoromethoxy)phenyl | 401 |
| 2,6-dimethylphenyl | cyclohexyl | 323 |
| 2,6-dimethylphenyl | 2-methylcyclohexyl | 337 |
| 2,6-dimethylphenyl | cycloheptyl | 667 |
| 2,6-dimethylphenyl | indan-1-yl | 357 |
| 2,6-dimethylphenyl | 2-dicyclohexyl | 406 |
| 2-phenoxyphenyl | phenyl | 381 |
| 2-phenoxyphenyl | o-tolyl | 395 |
| 2-phenoxyphenyl | 4-chloro-2-methylphenyl | 430 |
| 2-phenoxyphenyl | 2-fluorophenyl | 399 |
| 2-phenoxyphenyl | 3-fluorophenyl | 399 |
| 2-phenoxyphenyl | 4-fluorophenyl | 399 |
| 2-phenoxyphenyl | 2,4-difluorophenyl | 417 |
| 2-phenoxyphenyl | 2-methoxyphenyl | 411 |
| 2-phenoxyphenyl | 3-methoxyphenyl | 411 |
| 2-phenoxyphenyl | 4-methoxyphenyl | 411 |
| 2-phenoxyphenyl | 2,4-dimethoxyphenyl | 441 |
| 2-phenoxyphenyl | 2-phenoxyphenyl | 473 |
| 2-phenoxyphenyl | 3-phenoxyphenyl | 473 |
| 2-phenoxyphenyl | 4-phenoxyphenyl | 473 |
| 2-phenoxyphenyl | 4-biphenyl | 457 |
| 2-phenoxyphenyl | 4-benzylphenyl | 472 |
| 2-phenoxyphenyl | 4-(trifluoromethoxy)phenyl | 465 |
| 2-phenoxyphenyl | cyclohexyl | 387 |
| 2-phenoxyphenyl | 2-methylcyclohexyl | 401 |
| 2-phenoxyphenyl | cycloheptyl | 401 |
| 2-phenoxyphenyl | indan-1-yl | 421 |
| 2-phenoxyphenyl | 2-dicyclohexyl | 470 |

| | | |
|--------------------|----------------------------|-----|
| phenyl | phenyl | 289 |
| phenyl | o-tolyl | 303 |
| phenyl | 4-chloro-2-methylphenyl | 338 |
| phenyl | 2-fluorophenyl | 307 |
| phenyl | 3-fluorophenyl | 307 |
| phenyl | 4-fluorophenyl | 307 |
| phenyl | 2,4-difluorophenyl | 325 |
| phenyl | 2-methoxyphenyl | 319 |
| phenyl | 3-methoxyphenyl | 319 |
| phenyl | 4-methoxyphenyl | 319 |
| phenyl | 2,4-dimethoxyphenyl | 349 |
| phenyl | 2-phenoxyphenyl | 381 |
| phenyl | 3-phenoxyphenyl | 381 |
| phenyl | 4-phenoxyphenyl | 381 |
| phenyl | 4-biphenyl | 365 |
| phenyl | 4-benzylphenyl | 379 |
| phenyl | 4-(trifluoromethoxy)phenyl | 373 |
| phenyl | cyclohexyl | 295 |
| phenyl | 2-methylcyclohexyl | 309 |
| phenyl | cycloheptyl | 309 |
| phenyl | indan-1-yl | 329 |
| phenyl | 2-dicyclohexyl | 377 |
| 2,4-difluorophenyl | phenyl | 325 |
| 2,4-difluorophenyl | o-tolyl | 339 |
| 2,4-difluorophenyl | 4-chloro-2-methylphenyl | 374 |
| 2,4-difluorophenyl | 2-fluorophenyl | 343 |
| 2,4-difluorophenyl | 3-fluorophenyl | 343 |
| 2,4-difluorophenyl | 4-fluorophenyl | 343 |

| | | |
|-----------------------|----------------------------|-----|
| 2,4-difluorophenyl | 2,4-difluorophenyl | 361 |
| 2,4-difluorophenyl | 2-methoxyphenyl | 355 |
| 2,4-difluorophenyl | 3-methoxyphenyl | 355 |
| 2,4-difluorophenyl | 4-methoxyphenyl | 355 |
| 2,4-difluorophenyl | 2,4-dimethoxyphenyl | 385 |
| 2,4-difluorophenyl | 2-phenoxyphenyl | 417 |
| 2,4-difluorophenyl | 3-phenoxyphenyl | 417 |
| 2,4-difluorophenyl | 4-phenoxyphenyl | 417 |
| 2,4-difluorophenyl | 4-biphenyl | 401 |
| 2,4-difluorophenyl | 4-benzylphenyl | 415 |
| 2,4-difluorophenyl | 4-(trifluoromethoxy)phenyl | 409 |
| 2,4-difluorophenyl | cyclohexyl | 331 |
| 2,4-difluorophenyl | 2-methylcyclohexyl | 345 |
| 2,4-difluorophenyl | cycloheptyl | 345 |
| 2,4-difluorophenyl | indan-1-yl | 365 |
| 2,4-difluorophenyl | 2-dicyclohexyl | 413 |
| 2,6-diisopropylphenyl | phenyl | 373 |
| 2,6-diisopropylphenyl | o-tolyl | 387 |
| 2,6-diisopropylphenyl | 4-chloro-2-methylphenyl | 422 |
| 2,6-diisopropylphenyl | 2-fluorophenyl | 391 |
| 2,6-diisopropylphenyl | 3-fluorophenyl | 391 |
| 2,6-diisopropylphenyl | 4-fluorophenyl | 391 |
| 2,6-diisopropylphenyl | 2,4-difluorophenyl | 409 |
| 2,6-diisopropylphenyl | 2-methoxyphenyl | 403 |
| 2,6-diisopropylphenyl | 3-methoxyphenyl | 403 |
| 2,6-diisopropylphenyl | 4-methoxyphenyl | 403 |
| 2,6-diisopropylphenyl | 2,4-dimethoxyphenyl | 434 |
| 2,6-diisopropylphenyl | 2-phenoxyphenyl | 466 |

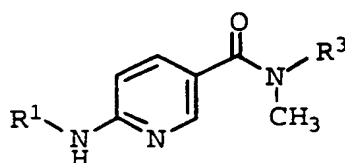
| | | |
|-----------------------|----------------------------|-----|
| 2,6-diisopropylphenyl | 3-phenoxyphenyl | 466 |
| 2,6-diisopropylphenyl | 4-phenoxyphenyl | 466 |
| 2,6-diisopropylphenyl | 4-biphenyl | 450 |
| 2,6-diisopropylphenyl | 4-benzylphenyl | 464 |
| 2,6-diisopropylphenyl | 4-(trifluoromethoxy)phenyl | 457 |
| 2,6-diisopropylphenyl | cyclohexyl | 380 |
| 2,6-diisopropylphenyl | 2-methylcyclohexyl | 394 |
| 2,6-diisopropylphenyl | cycloheptyl | 394 |
| 2,6-diisopropylphenyl | indan-1-yl | 414 |
| 2,6-diisopropylphenyl | 2-dicyclohexyl | 462 |
| 2,4-dimethoxyphenyl | phenyl | 349 |
| 2,4-dimethoxyphenyl | o-tolyl | 363 |
| 2,4-dimethoxyphenyl | 4-chloro-2-methylphenyl | 398 |
| 2,4-dimethoxyphenyl | 2-fluorophenyl | 367 |
| 2,4-dimethoxyphenyl | 3-fluorophenyl | 367 |
| 2,4-dimethoxyphenyl | 4-fluorophenyl | 367 |
| 2,4-dimethoxyphenyl | 2,4-difluorophenyl | 385 |
| 2,4-dimethoxyphenyl | 2-methoxyphenyl | 379 |
| 2,4-dimethoxyphenyl | 3-methoxyphenyl | 379 |
| 2,4-dimethoxyphenyl | 4-methoxyphenyl | 379 |
| 2,4-dimethoxyphenyl | 2,4-dimethoxyphenyl | 409 |
| 2,4-dimethoxyphenyl | 2-phenoxyphenyl | 441 |
| 2,4-dimethoxyphenyl | 3-phenoxyphenyl | 441 |
| 2,4-dimethoxyphenyl | 4-phenoxyphenyl | 441 |
| 2,4-dimethoxyphenyl | 4-biphenyl | 425 |
| 2,4-dimethoxyphenyl | 4-benzylphenyl | 439 |
| 2,4-dimethoxyphenyl | 4-(trifluoromethoxy)phenyl | 433 |
| 2,4-dimethoxyphenyl | 3-trifluoromethylphenyl | 417 |

| | | |
|---------------------|----------------------------|-----|
| 2,4-dimethoxyphenyl | cyclohexyl | 355 |
| 2,4-dimethoxyphenyl | 2-methylcyclohexyl | 369 |
| 3-methoxyphenyl | phenyl | 319 |
| 3-methoxyphenyl | o-tolyl | 333 |
| 3-methoxyphenyl | 4-chloro-2-methylphenyl | 368 |
| 3-methoxyphenyl | 2-fluorophenyl | 337 |
| 3-methoxyphenyl | 3-fluorophenyl | 337 |
| 3-methoxyphenyl | 4-fluorophenyl | 337 |
| 3-methoxyphenyl | 2,4-difluorophenyl | 355 |
| 3-methoxyphenyl | 2-methoxyphenyl | 349 |
| 3-methoxyphenyl | 3-methoxyphenyl | 349 |
| 3-methoxyphenyl | 4-methoxyphenyl | 349 |
| 3-methoxyphenyl | 2,4-dimethoxyphenyl | 379 |
| 3-methoxyphenyl | 2-phenoxyphenyl | 411 |
| 3-methoxyphenyl | 3-phenoxyphenyl | 411 |
| 3-methoxyphenyl | 4-phenoxyphenyl | 411 |
| 3-methoxyphenyl | 4-biphenyl | 395 |
| 3-methoxyphenyl | 4-benzylphenyl | 409 |
| 3-methoxyphenyl | 4-(trifluoromethoxy)phenyl | 403 |
| 3-methoxyphenyl | 3-trifluoromethylphenyl | 387 |
| 3-methoxyphenyl | cyclohexyl | 625 |
| 3-methoxyphenyl | 2-methylcyclohexyl | 339 |
| 4-methoxyphenyl | phenyl | 319 |
| 4-methoxyphenyl | o-tolyl | 333 |
| 4-methoxyphenyl | 4-chloro-2-methylphenyl | 368 |
| 4-methoxyphenyl | 2-fluorophenyl | 337 |
| 4-methoxyphenyl | 3-fluorophenyl | 337 |
| 4-methoxyphenyl | 4-fluorophenyl | 337 |

| | | |
|-----------------|----------------------------|-----|
| 4-methoxyphenyl | 2,4-difluorophenyl | 355 |
| 4-methoxyphenyl | 2-methoxyphenyl | 349 |
| 4-methoxyphenyl | 3-methoxyphenyl | 349 |
| 4-methoxyphenyl | 4-methoxyphenyl | 349 |
| 4-methoxyphenyl | 2,4-dimethoxyphenyl | 379 |
| 4-methoxyphenyl | 2-phenoxyphenyl | 411 |
| 4-methoxyphenyl | 3-phenoxyphenyl | 411 |
| 4-methoxyphenyl | 4-phenoxyphenyl | 411 |
| 4-methoxyphenyl | 4-biphenyl | 395 |
| 4-methoxyphenyl | 4-benzylphenyl | 409 |
| 4-methoxyphenyl | 4-(trifluoromethoxy)phenyl | 403 |
| 4-methoxyphenyl | 3-trifluoromethylphenyl | 387 |
| 4-methoxyphenyl | cyclohexyl | 625 |
| 4-methoxyphenyl | 2-methylcyclohexyl | 339 |
| 2-methoxyphenyl | phenyl | 319 |
| 2-methoxyphenyl | o-tolyl | 333 |
| 2-methoxyphenyl | 4-chloro-2-methylphenyl | 368 |
| 2-methoxyphenyl | 2-fluorophenyl | 337 |
| 2-methoxyphenyl | 3-fluorophenyl | 337 |
| 2-methoxyphenyl | 4-fluorophenyl | 337 |
| 2-methoxyphenyl | 2,4-difluorophenyl | 355 |
| 2-methoxyphenyl | 2-methoxyphenyl | 349 |
| 2-methoxyphenyl | 3-methoxyphenyl | 349 |
| 2-methoxyphenyl | 4-methoxyphenyl | 349 |
| 2-methoxyphenyl | 2,4-dimethoxyphenyl | 379 |
| 2-methoxyphenyl | 2-phenoxyphenyl | 411 |
| 2-methoxyphenyl | 3-phenoxyphenyl | 411 |
| 2-methoxyphenyl | 4-phenoxyphenyl | 411 |

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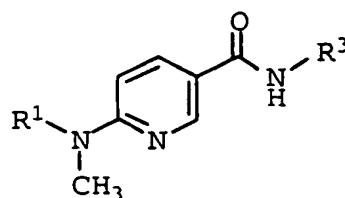
| | | |
|-----------------|----------------------------|-----|
| 2-methoxyphenyl | 4-biphenyl | 395 |
| 2-methoxyphenyl | 4-benzylphenyl | 409 |
| 2-methoxyphenyl | 4-(trifluoromethoxy)phenyl | 403 |
| 2-methoxyphenyl | 3-trifluoromethylphenyl | 387 |
| 2-methoxyphenyl | cyclohexyl | 625 |
| 2-methoxyphenyl | 2-methylcyclohexyl | 339 |

Table 9

| <u>R³</u> | <u>R¹</u> | <u>MS</u> <u>(m/z)</u> |
|----------------------|-------------------------|---------------------------|
| 4-chlorophenyl | phenyl | 338 |
| 4-chlorophenyl | o-tolyl | 352 |
| 4-chlorophenyl | 4-chloro-2-methylphenyl | 386 |
| 4-chlorophenyl | 2-fluorophenyl | 356 |
| 4-chlorophenyl | 3-fluorophenyl | 356 |
| 4-chlorophenyl | 4-fluorophenyl | 356 |
| 4-chlorophenyl | 2,4-difluorophenyl | 374 |
| 4-chlorophenyl | 2-methoxyphenyl | 368 |
| 4-chlorophenyl | 3-methoxyphenyl | 368 |
| 4-chlorophenyl | 4-methoxyphenyl | 368 |
| 4-chlorophenyl | 2,4-dimethoxyphenyl | 398 |
| 4-chlorophenyl | 2-phenoxyphenyl | 430 |
| 4-chlorophenyl | 3-phenoxyphenyl | 430 |
| 4-chlorophenyl | 4-phenoxyphenyl | 430 |
| 4-chlorophenyl | 4-biphenyl | 414 |

| | | |
|----------------|----------------------------|-----|
| 4-chlorophenyl | 4-benzylphenyl | 428 |
| 4-chlorophenyl | 4-(trifluoromethoxy)phenyl | 422 |
| 4-chlorophenyl | cyclohexyl | 344 |
| 4-chlorophenyl | 2-methylcyclohexyl | 358 |
| phenyl | phenyl | 303 |
| phenyl | o-tolyl | 317 |
| phenyl | 4-chloro-2-methylphenyl | 352 |
| phenyl | 2-fluorophenyl | 321 |
| phenyl | 3-fluorophenyl | 321 |
| phenyl | 4-fluorophenyl | 321 |
| phenyl | 2,4-difluorophenyl | 339 |
| phenyl | 2-methoxyphenyl | 333 |
| phenyl | 3-methoxyphenyl | 333 |
| phenyl | 4-methoxyphenyl | 333 |
| phenyl | 2,4-dimethoxyphenyl | 363 |
| phenyl | 2-phenoxyphenyl | 395 |
| phenyl | 3-phenoxyphenyl | 395 |
| phenyl | 4-phenoxyphenyl | 395 |
| phenyl | 4-biphenyl | 379 |
| phenyl | 4-benzylphenyl | 393 |
| phenyl | 4-(trifluoromethoxy)phenyl | 387 |
| phenyl | 3-trifluoromethylphenyl | 371 |
| phenyl | cyclohexyl | 309 |
| phenyl | 2-methylcyclohexyl | 323 |

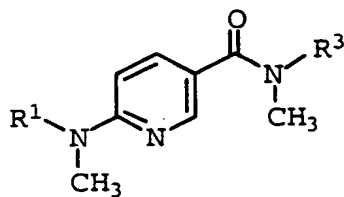
Table 10



| <u>R³</u> | <u>R¹</u> | <u>MS</u> <u>(m/z)</u> |
|-----------------------|-------------------------|---------------------------|
| o-tolyl | N-methylphenyl | 317 |
| o-tolyl | 4-chloro-N-methylphenyl | 352 |
| o-tolyl | N-methylcyclohexyl | 323 |
| 2-fluorophenyl | N-methylphenyl | 321 |
| 2-fluorophenyl | 4-chloro-N-methylphenyl | 356 |
| 2-fluorophenyl | N-methylcyclohexyl | 327 |
| 2,6-dimethylphenyl | N-methylphenyl | 331 |
| 2,6-dimethylphenyl | 4-chloro-N-methylphenyl | 366 |
| 2,6-dimethylphenyl | N-methylcyclohexyl | 337 |
| 2-phenoxyphenyl | N-methylphenyl | 395 |
| 2-phenoxyphenyl | 4-chloro-N-methylphenyl | 430 |
| 2-phenoxyphenyl | N-methylcyclohexyl | 401 |
| phenyl | N-methylphenyl | 303 |
| phenyl | 4-chloro-N-methylphenyl | 338 |
| phenyl | N-methylcyclohexyl | 309 |
| 2,4-difluorophenyl | N-methylphenyl | 339 |
| 2,4-difluorophenyl | 4-chloro-N-methylphenyl | 374 |
| 2,4-difluorophenyl | N-methylcyclohexyl | 345 |
| 2,6-diisopropylphenyl | N-methylphenyl | 387 |
| 2,6-diisopropylphenyl | 4-chloro-N-methylphenyl | 422 |
| 2,6-diisopropylphenyl | N-methylcyclohexyl | 394 |

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| | | |
|---------------------|-------------------------|-----|
| 2,4-dimethoxyphenyl | N-methylphenyl | 363 |
| 2,4-dimethoxyphenyl | 4-chloro-N-methylphenyl | 398 |
| 2,4-dimethoxyphenyl | N-methylcyclohexyl | 369 |
| 3-methoxyphenyl | N-methylphenyl | 333 |
| 3-methoxyphenyl | 4-chloro-N-methylphenyl | 368 |
| 3-methoxyphenyl | N-methylcyclohexyl | 339 |
| 4-methoxyphenyl | N-methylphenyl | 333 |
| 4-methoxyphenyl | 4-chloro-N-methylphenyl | 368 |
| 4-methoxyphenyl | N-methylcyclohexyl | 339 |
| 2-methoxyphenyl | N-methylphenyl | 333 |
| 2-methoxyphenyl | 4-chloro-N-methylphenyl | 368 |
| 2-methoxyphenyl | N-methylcyclohexyl | 339 |

Table 11

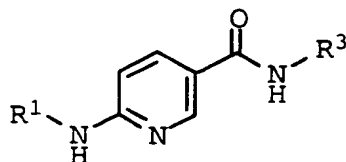
| <u>R¹</u> | <u>R¹</u> | <u>MS</u> <u>(m/z)</u> |
|----------------------|----------------------|---------------------------|
| 4-chlorophenyl | phenyl | 352 |
| 4-chlorophenyl | 4-chlorophenyl | 386 |
| 4-chlorophenyl | cyclohexyl | 358 |
| phenyl | phenyl | 317 |
| phenyl | 4-chlorophenyl | 352 |
| phenyl | cyclohexyl | 323 |

5

The compounds listed in Tables 12-13 can be prepared from 6-chloro-N-substituted nicotinamides

compounds and the appropriate amine according to the general procedure above.

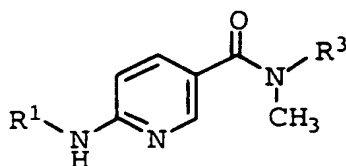
Table 12



5

| <u>R¹X</u> | <u>R³</u> |
|-----------------------|-------------------------|
| o-tolyl | 3-trifluoromethylphenyl |
| 2-fluorophenyl | 3-trifluoromethylphenyl |
| 2,6-dimethylphenyl | 3-trifluoromethylphenyl |
| 2-phenoxyphenyl | 3-trifluoromethylphenyl |
| phenyl | 3-trifluoromethylphenyl |
| 2,4-difluorophenyl | 3-trifluoromethylphenyl |
| 2,6-diisopropylphenyl | 3-trifluoromethylphenyl |
| 2,4-dimethoxyphenyl | cycloheptyl |
| 2,4-dimethoxyphenyl | indan-1-yl |
| 2,4-dimethoxyphenyl | 2-dicyclohexyl |
| 3-methoxyphenyl | cycloheptyl |
| 3-methoxyphenyl | indan-1-yl |
| 3-methoxyphenyl | 2-dicyclohexyl |
| 4-methoxyphenyl | cycloheptyl |
| 4-methoxyphenyl | indan-1-yl |
| 4-methoxyphenyl | 2-dicyclohexyl |
| 2-methoxyphenyl | cycloheptyl |
| 2-methoxyphenyl | indan-1-yl |
| 2-methoxyphenyl | 2-dicyclohexyl |

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Table 13

| <u>R¹X</u> | <u>R³</u> |
|-----------------------|-------------------------|
| 4-chlorophenyl | 3-trifluoromethylphenyl |
| 4-chlorophenyl | cycloheptyl |
| 4-chlorophenyl | indan-1-yl |
| 4-chlorophenyl | 2-dicyclohexyl |
| phenyl | cycloheptyl |
| phenyl | indan-1-yl |
| phenyl | 2-dicyclohexyl |

5

Example 20

The following assays were used to characterize the ability of compounds of the invention to inhibit the production of TNF- α and IL-1- β . The second assay

10 measured the inhibition of TNF- α and/or IL-1- β in mice after oral administration of the test compounds. The third assay, a glucagon binding inhibition in vitro assay, can be used to characterize the ability of compounds of the invention to inhibit glucagon binding.

15 The fourth assay, a Cyclooxygenase enzyme (COX-1 and COX-2) inhibition activity in vitro assay, can be used to characterize the ability of compounds of the invention to inhibit COX-1 and/or COX-2. The fifth

20 assay, a Raf-kinase inhibition assay, can be used to characterize the compounds of the invention to inhibit phosphorylation of MEK by activated Raf-kinase.

Lipopolysaccharide-activated monocyte TNF production assay

Isolation of monocytes

5 Test compounds were evaluated *in vitro* for the ability to inhibit the production of tumor necrosis factor (TNF) by monocytes activated with bacterial lipopolysaccharide (LPS). Fresh residual source leukocytes (a byproduct of plateletpheresis) were
10 obtained from a local blood bank, and peripheral blood mononuclear cells (PBMCs) were isolated by density gradient centrifugation on Ficol-Paque Plus (Pharmacia). PBMCs were suspended at 2×10^6 /ml in DMEM supplemented to contain 2% FCS, 10 mM, 0.3 mg/ml glutamate, 100 U/ml
15 penicillin G and 100 mg/ml streptomycin sulfate (complete media). Cells were plated into Falcon flat bottom, 96 well culture plates (200 μ l/well) and cultured overnight at 37°C and 6% CO₂. Non-adherent cells were removed by washing with 200 μ l/well of fresh
20 medium. Wells containing adherent cells (~70% monocytes) were replenished with 100 μ l of fresh medium.

Preparation of test compound stock solutions

 Test compounds were dissolved in DMSO. Compound
25 stock solutions were prepared to an initial concentration of 10 - 50 μ M. Stocks were diluted initially to 20 - 200 μ M in complete media. Nine two-fold serial dilutions of each compound were then prepared in complete medium.

30

Treatment of cells with test compounds and activation of TNF production with lipopolysaccharide

 One hundred microliters of each test compound dilution were added to microtiter wells containing
35 adherent monocytes and 100 μ l complete medium. Monocytes were cultured with test compounds for 60 min at which time 25 μ l of complete medium containing 30

ng/ml lipopolysaccharide from *E. coli* K532 were added to each well. Cells were cultured an additional 4 hrs. Culture supernatants were then removed and TNF presence in the supernatants was quantified using an ELISA.

5

TNF ELISA

Flat bottom, 96 well Corning High Binding ELISA plates were coated overnight (4°C) with 150 µL/well of 3 µg/ml murine anti-human TNF-α MAb (R&D Systems #MAB210).
10 Wells were then blocked for 1 hr at room temperature with 200 µL/well of CaCl₂-free ELISA buffer supplemented to contain 20 mg/ml BSA (standard ELISA buffer: 20 mM, 150 mM NaCl, 2 mM CaCl₂, 0.15 mM thimerosal, pH 7.4). Plates were washed and replenished with 100 µl of test
15 supernatants (diluted 1:3) or standards. Standards consisted of eleven 1.5-fold serial dilutions from a stock of 1 ng/ml recombinant human TNF (R&D Systems). Plates were incubated at room temperature for 1 hr on orbital shaker (300 rpm), washed and replenished with
20 100 µl/well of 0.5 µg/ml goat anti-human TNF-α (R&D systems #AB-210-NA) biotinylated at a 4:1 ratio. Plates were incubated for 40 min, washed and replenished with 100 µl/well of alkaline phosphatase-conjugated streptavidin (Jackson ImmunoResearch #016-050-084) at
25 0.02 µg/ml. Plates were incubated 30 min, washed and replenished with 200 µl/well of 1 mg/ml of p-nitrophenyl phosphate. After 30 min, plates were read at 405 nm on a V_{max} plate reader.

30 Data analysis

Standard curve data were fit to a second order polynomial and unknown TNF-α concentrations determined from their OD by solving this equation for concentration. TNF concentrations were then plotted vs.
35 test compound concentration using a second order polynomial. This equation was then used to calculate

the concentration of test compounds causing a 50% reduction in TNF production.

- The following compounds had an IC_{50} of less than 15 μM :
- 5 2-cyclohexyloxy-5-(2-chlorophenylcarbonylamino)pyridine;
 - 2-cyclohexyloxy-5-(2-methylphenylcarbonylamino)pyridine;
 - 2-cyclohexyloxy-5-(2,6-dichlorophenylcarbonylamino)pyridine;
 - 10 2-cyclohexyloxy-5-(2,6-dimethylphenylcarbonylamino)pyridine;
 - 2-(2,4-dimethylphenoxy)-5-(2-methylphenylcarbonylamino)pyridine;
 - 2-(2-methyl-4-fluorophenoxy)-5-(2-methylphenylcarbonyl
 - 15 amino)pyridine;
 - 2-(2-methyl-4-chlorophenoxy)-5-(2-chlorophenylcarbonyl amino)pyridine;
 - 2-(2-methyl-4-chlorophenoxy)-5-(2,6-dichlorophenyl carbonylamino)pyridine;
 - 20 2-(2-methyl-4-chlorophenoxy)-5-(2,6-dimethylphenyl carbonylamino)pyridine;
 - 2-(4-chlorophenoxy)-5-(2,6-dimethylphenylcarbonylamino)pyridine;
 - 2-(2-methyl-4-fluorophenoxy)-5-(2,6-dichlorophenyl
 - 25 carbonylamino)pyridine;
 - 2-(2-methyl-4-fluorophenoxy)-5-(2,6-dimethylphenyl carbonylamino)pyridine;
 - 2-(2-methyl-4-fluorophenoxy)-5-(2-fluorophenylcarbonyl amino)pyridine;
 - 30 2-(2,4-dimethylphenoxy)-5-(2,6-dimethylphenylcarbonyl amino)pyridine;
 - 2-(1-naphthyloxy)-5-(2-methylphenylcarbonylamino)pyridine;
 - 2-(1-naphthyloxy)-5-(2,6-dichlorophenylcarbonylamino)
 - 35 pyridine;
 - 2-(2-methyl-3-pyridyloxy)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;

- 2-(2-methyl-4-chlorophenoxy)-5-((3,5-dimethyl-4-isoxazolyl)carbonylamino)pyridine;
2-cyclohexylamino-5-(2,6-dichlorophenylcarbonylamino)pyridine;
5 2-cyclohexylamino-5-(2,6-dimethylphenylcarbonylamino)pyridine;
2-(2-methylcyclohexylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
2-(2-methylcyclohexylamino)-5-(2-methylphenylcarbonyl
10 amino)pyridine;
2-(2-methylphenylamino)-5-(2-methylphenylcarbonyl amino)pyridine;
2-(2-methylphenylamino)-5-(2,6-dimethylphenylcarbonyl amino)pyridine;
15 2-(2-methyl-4-chlorophenylamino)-5-(2-methylphenyl carbonylamino)pyridine; and
2-(2-methyl-4-chlorophenylamino)-5-(2-methylphenyl aminocarbonyl)pyridine.
- 20 Compounds of the invention can also be shown to inhibit LPS-induced release of IL-1 β , IL-6 and/or IL-8 from monocytes by measuring concentrations of IL-1 β , IL-6 and/or IL-8 by methods well known to those skilled in the art. In a similar manner to the above described
25 assay involving the LPS induced release of TNF- α from monocytes, compounds of this invention can also be shown to inhibit LPS induced release of IL-1 β , IL-6 and/or IL-8 from monocytes by measuring concentrations of IL-1 β , IL-6 and/or IL-8 by methods well known to those skilled
30 in the art. Thus, the compounds of the invention may lower elevated levels of TNF- α , IL-1, IL-6, and IL-8 levels. Reducing elevated levels of these inflammatory cytokines to basal levels or below is favorable in controlling, slowing progression, and alleviating many
35 disease states. All of the compounds are useful in the methods of treating disease states in which TNF- α , IL-

1 β , IL-6, and IL-8 play a role to the full extent of the definition of TNF- α -mediated diseases described herein.

Inhibition of LPS-Induced TNF- α production in mice

5 Male DBA/1LACJ mice are dosed with vehicle or test compounds in a vehicle (the vehicle consisting of 0.5% tragacanth in 0.03 N HCl) 30 minutes prior to lipopolysaccharide (2 mg/kg, I.V.) injection. Ninety minutes after LPS injection, blood are collected and the
10 serum is analyzed by ELISA for TNF levels.

 Selected compounds from the class have shown in vivo activity in a LPS mouse model in which serum levels of TNF- α were reduced in the presence of compounds of this invention.

15 Compounds of the invention may be shown to have anti-inflammatory properties in animal models of inflammation, including carageenan paw edema, collagen induced arthritis and adjuvant arthritis, such as the carageenan paw edema model (C. A. Winter et al Proc.
20 Soc. Exp. Biol. Med. (1962) vol 111, p 544; K. F. Swingle, in R. A. Scherrer and M. W. Whitehouse, Eds., Antiinflammatory Agents, Chemistry and Pharmacology, Vol. 13-II, Academic, New York, 1974, p. 33) and collagen induced arthritis (D. E. Trentham et al J. Exp.
25 Med. (1977) vol. 146, p 857; J. S. Courtenay, Nature (New Biol.) (1980), Vol 283, p 666).

¹²⁵I-Glucagon Binding Screen with CHO/hGLUR Cells

 The assay is described in WO 97/16442, which is
30 incorporated herein by reference in its entirety.

Reagents

 The reagents can be prepared as follows: (a) prepare fresh 1M o-Phenanthroline (Aldrich) (198.2 mg/ml
35 ethanol); (b) prepare fresh 0.5M DTT (Sigma); (c) Protease Inhibitor Mix (1000X): 5 mg leupeptin, 10 mg

benzamidine, 40 mg bacitracin and 5 mg soybean trypsin inhibitor per ml DMSO and store aliquots at -20°C; (d) 250 µM human glucagon (Peninsula): solubilize 0.5 mg vial in 575 µl 0.1N acetic acid (1 µl yields 1 µM final concentration in assay for non-specific binding) and store in aliquots at -20°C; (e) Assay Buffer: 20mM Tris (pH 7.8), 1 mM DTT and 3 mM o-phenanthroline; (f) Assay Buffer with 0.1% BSA (for dilution of label only; 0.01% final in assay): 10 µl 10% BSA (heat-inactivated) and 990 µl Assay Buffer; (g) ¹²⁵I-Glucagon (NEN, receptor-grade, 2200 Ci/mmol): dilute to 50,000 cpm/25 µl in assay buffer with BSA (about 50pM final concentration in assay).

15 Harvesting of CHO/hGLUR Cells for Assay

1. Remove media from confluent flask then rinse once each with PBS (Ca, Mg-free) and Enzyme-free Dissociation Fluid (Specialty Media, Inc.).
2. Add 10 ml Enzyme-free Dissoc. Fluid and hold for about 4 min. at 37°C.
3. Gently tap cells free, triturate, take aliquot for counting and centrifuge remainder for 5 min. at 1000 rpm.
4. Resuspend pellet in Assay Buffer at 75000 cells per 100 µl.

Membrane preparations of CHO/hGLUR cells can be used in place of whole cells at the same assay volume. Final protein concentration of a membrane preparation is determined on a per batch basis.

30

Assay

- The determination of inhibition of glucagon binding can be carried out by measuring the reduction of I¹²⁵-glucagon binding in the presence of compounds of Formula I. The reagents are combined in 120 µL of assay buffer as follows:

| | Compound/ Vehicle | 250 μ M Glucagon | 125 I- Glucagon | CHO/hGLUR Cells |
|---------------------|----------------------|-------------------------|-------------------------|--------------------|
| Total Binding | --/5 μ l | -- | 25 μ l | 100 μ l |
| + Compound | 5 μ l/-- | -- | 25 μ l | 100 μ l |
| Nonspecific Binding | --/5 μ l | 1 μ l | 25 μ l | 100 μ l |

The mixture is incubated for 60 min. at 22°C on a shaker at 275 rpm. The mixture is filtered over pre-soaked (0.5% polyethylimine (PEI)) GF/C filtermat using an Innotech Harvester or Tomtec Harvester with four washes of ice-cold 20mM Tris buffer (pH 7.8). The radioactivity in the filters is determined by a gamma-scintillation counter.

Thus, compounds of the invention may also be shown to inhibit the binding of glucagon to glucagon receptors.

Cyclooxygenase Enzyme Activity Assay

The human monocytic leukemia cell line, THP-1, differentiated by exposure to phorbol esters expresses only COX-1; the human osteosarcoma cell line 143B expresses predominantly COX-2. THP-1 cells are routinely cultured in RPMI complete media supplemented with 10% FBS and human osteosarcoma cells (HOSC) are cultured in minimal essential media supplemented with 10% fetal bovine serum (MEM-10%FBS); all cell incubations are at 37°C in a humidified environment containing 5% CO₂.

COX-1 Assay

In preparation for the COX-1 assay, THP-1 cells are grown to confluency, split 1:3 into RPMI containing 2%

FBS and 10 mM phorbol 12-myristate 13-acetate (TPA), and incubated for 48 hours on a shaker to prevent attachment. Cells are pelleted and resuspended in Hank's Buffered Saline (HBS) at a concentration of 2.5×10^6 cells/mL and plated in 96-well culture plates at a density of 5×10^5 cells/mL. Test compounds are diluted in HBS and added to the desired final concentration and the cells are incubated for an additional 4 hours. Arachidonic acid is added to a final concentration of 30 mM, the cells incubated for 20 minutes at 37°C, and enzyme activity determined as described below.

COX-2 Assay

For the COX-2 assay, subconfluent HOSC are trypsinized and resuspended at 3×10^6 cells/mL in MEM-FBS containing 1 ng human IL-1b/mL, plated in 96-well tissue culture plates at a density of 3×10^4 cells per well, incubated on a shaker for 1 hour to evenly distribute cells, followed by an additional 2 hour static incubation to allow attachment. The media is then replaced with MEM containing 2% FBS (MEM-2%FBS) and 1 ng human IL-1b/mL, and the cells incubated for 18-22 hours. Following replacement of media with 190 mL MEM, 10 mL of test compound diluted in HBS is added to achieve the desired concentration and the cells incubated for 4 hours. The supernatants are removed and replaced with MEM containing 30 mM arachidonic acid, the cells incubated for 20 minutes at 37°C, and enzyme activity determined as described below.

COX Activity Determined

After incubation with arachidonic acid, the reactions are stopped by the addition of 1 N HCl, followed by neutralization with 1 N NaOH and centrifugation to pellet cell debris. Cyclooxygenase enzyme activity in both HOSC and THP-1 cell supernatants

is determined by measuring the concentration of PGE₂ using a commercially available ELISA (Neogen #404110). A standard curve of PGE₂ is used for calibration, and commercially available COX-1 and COX-2 inhibitors are included as standard controls.

The following compound exhibits activities in the Cyclooxygenase assay with IC₅₀ values of 10 μ M or less: 2-(2,4-dimethylphenylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine.

10

Raf Kinase assay

In vitro Raf kinase activity is measured by the extent of phosphorylation of the substrate MEK (Map kinase/ERK kinase) by activated Raf kinase.

15 Phosphorylated MEK is trapped on a filter and incorporation of radiolabeled phosphate is quantified by scintillation counting.

MATERIALS:

20

Activated Raf is produced by triple transfection of Sf9 cells with baculoviruses expressing "Glu-Glu"-epitope tagged Raf, val¹²-H-Ras, and Lck. The "Glu-Glu"-epitope, Glu-Try-Met-Pro-Met-Glu, was fused to the carboxy-terminus of full length c-Raf.

25

Catalytically inactive MEK (K97A mutation) is produced in Sf9 cells transfected with a baculovirus expressing c-terminus "Glu-Glu" epitope-tagged K97A MEK1.

Anti "Glu-Glu" antibody was purified from cells grown as described in: Grussenmeyer, et al., Proceedings of the National Academy of Science, U.S.A. pp 7952-7954, 1985. Column buffer: 20 mM Tris pH=8, 100 mM NaCl, 1 mM EDTA, 2.5 mM EGTA, 10 mM MgCl₂, 2 mM DTT, 0.4 mM AEBSF, 0.1% n-octylglucopyranoside, 1 mM okadaic acid, and 10 μ g/mL each of benzamidine, leupeptin, pepstatin, and aprotinin.

35

5x Reaction buffer: 125 mM HEPES pH=8, 25 mM MgCl₂, 5 mM EDTA, 5 mM Na₂VO₄, 100 µg/mL BSA.

Enzyme dilution buffer: 25 mM HEPES pH=8, 1 mM EDTA, 1 mM Na₂VO₄, 400 µg/mL BSA.

- 5 Stop solution: 100 mM EDTA, 80 mM sodium pyrophosphate.
Filter plates: Milipore multiscreen # SE3MO78E3,
Immobilon-P (PVDF).

METHODS:

10

Protein purification: Sf9 cells were infected with baculovirus and grown as described in Williams, et al., Proceedings of the National Academy of Science, U.S.A. pp 2922-2926, 1992. All subsequent steps were preformed
15 on ice or at 4 °C. Cells were pelleted and lysed by sonication in column buffer. Lysates were spun at 17,000xg for 20 min, followed by 0.22 µm filtration. Epitope tagged proteins were purified by chromatography over GammaBind Plus affinity column to which the "Glu-Glu"
20 antibody was coupled. Proteins were loaded on the column followed by sequential washes with two column volumes of column buffer, and eluted with 50 µg/mL Glu-Tyr-Met-Pro-Met-Glu in column buffer.

- 25 Raf kinase assay: Test compounds were evaluated using ten 3-fold serial dilutions starting at 10 - 100 µM. 10 µL of the test inhibitor or control, dissolved in 10% DMSO, was added to the assay plate followed by the addition of 30 µL of the a mixture containing 10 µL 5x
30 reaction buffer, 1mM ³³P-γ-ATP (20 µCi/mL), 0.5 µL MEK (2.5 mg/mL), 1 µL 50 mM β-mercaptoethanol. The reaction was started by the addition of 10 µL of enzyme dilution buffer containing 1 mM DTT and an amount of activated Raf that produces linear kinetics over the reaction time
35 course. The reaction was mixed and incubated at room

temperature for 90 min. and stopped by the addition of 50 μ L stop solution. 90 μ L aliquots of this stopped solution were transferred onto GFP-30 cellulose microtiter filter plates (Polyfiltronics), the filter plates washed in four well volumes of 5% phosphoric acid, allowed to dry, and then replenished with 25 μ L scintillation cocktail. The plates were counted for ^{33}P gamma emission using a TopCount Scintillation Reader.

Accordingly, the compounds of the invention or a pharmaceutical composition thereof are useful for prophylaxis and treatment of rheumatoid arthritis; Pagets disease; osteoporosis; multiple myeloma; uveitis; acute and chronic myelogenous leukemia; pancreatic β cell destruction; osteoarthritis; rheumatoid spondylitis; gouty arthritis; inflammatory bowel disease; adult respiratory distress syndrome (ARDS); psoriasis; Crohn's disease; allergic rhinitis; ulcerative colitis; anaphylaxis; contact dermatitis; asthma; muscle degeneration; cachexia; Reiter's syndrome; type I and type II diabetes; bone resorption diseases; graft vs. host reaction; ischemia reperfusion injury; atherosclerosis; brain trauma; Alzheimer's disease; stroke; myocardial infarction; multiple sclerosis; cerebral malaria; sepsis; septic shock; toxic shock syndrome; fever, and myalgias due to infection. HIV-1, HIV-2, HIV-3, cytomegalovirus (CMV), influenza, adenovirus, the herpes viruses (including HSV-1, HSV-2), and herpes zoster, all of which are sensitive to TNF- α and/or IL-1 inhibition or glucagon antagonism, will also be positively effected by the compounds and methods of the invention.

The compounds of the present invention may also possess oncolytic characteristics and may be useful for the treatment of cancer. The compounds of the present invention may also block signal transduction by extracellular mitogenic stimuli and oncoproteins through inhibition of Raf kinase.

The compounds of the present invention also may possess analgesic properties and may be useful for the treatment of pain disorders, such as hyperalgesia due to excessive IL-1. The compounds of the present invention
5 may also prevent the production of prostaglandins by inhibition of enzymes in the human arachidonic acid/prostaglandin pathway, including cyclooxygenase (WO 96/03387, incorporated herein by reference in its entirety).

10 Because of their ability to lower TNF- α and IL-1 concentrations or inhibit glucagon binding to its receptor, the compounds of the invention are also useful research tools for studying the physiology associated with blocking these effects.

15 The methods of the invention comprise administering an effective dose of a compound of the invention, a pharmaceutical salt thereof, or a pharmaceutical composition of either, to a subject (i.e., an animal, preferably a mammal, most preferably a human) in need of
20 a reduction in the level of TNF- α , IL-1, IL-6, and/or IL-8 levels and/or reduction in plasma glucose levels and/or which subject may be suffering from rheumatoid arthritis; Pagets disease; osteoporosis; multiple myeloma; uveitis; acute and chronic myelogenous
25 leukemia; pancreatic β cell destruction; osteoarthritis; rheumatoid spondylitis; gouty arthritis; inflammatory bowel disease; adult respiratory distress syndrome (ARDS); psoriasis; Crohn's disease; allergic rhinitis; ulcerative colitis; anaphylaxis; contact dermatitis;
30 asthma; muscle degeneration; cachexia; Reiter's syndrome; type I and type II diabetes; cancer; bone resorption diseases; graft vs. host reaction; Alzheimer's disease; stroke; myocardial infarction; ischemia reperfusion injury; atherosclerosis; brain
35 trauma; multiple sclerosis; cerebral malaria; sepsis; septic shock; toxic shock syndrome; fever, and myalgias due to infection, or which subject is infected by HIV-1,

HIV-2, HIV-3, cytomegalovirus (CMV), influenza, adenovirus, the herpes viruses (including HSV-1, HSV-2), or herpes zoster.

In another aspect, this invention comprises the use
5 of a compound of the invention, or pharmaceutically acceptable salts thereof, in the manufacture of a medicament for the treatment either acutely or chronically of a TNF- α , IL-1 β , IL-6, and/or IL-8 mediated disease state, including those described
10 previously. The compounds of the present are also useful in the manufacture of an anti-cancer medicament. The compounds of the present invention are also useful in the manufacture of a medicament to attenuate or prevent signal transduction by extracellular mitogenic stimuli
15 and oncoproteins through inhibition of Raf kinase. Also, the compounds of this invention are useful in the manufacture of an analgesic medicament and a medicament for treating pain disorders, such as hyperalgesia. The compounds of the present invention also are useful in
20 the manufacture of a medicament to prevent the production of prostaglandins by inhibition of enzymes in the human arachidonic acid/prostaglandin pathway.

In still another aspect, this invention provides a pharmaceutical composition comprising an effective TNF-
25 α , IL-1 β , IL-6, and/or IL-8 lowering amount and/or effective plasma glucose level lowering amount, and/or effective tumor suppressing amount of a compound of the invention and a pharmaceutically acceptable carrier or diluent, and if desired other active ingredients. The
30 compounds of the invention are administered by any suitable route, preferably in the form of a pharmaceutical composition adapted to such a route, and in a dose effective for the treatment intended. Therapeutically effective doses of the compounds of the
35 present invention required to arrest the progress or prevent tissue damage associated with the disease are

readily ascertained by one of ordinary skill in the art using standard methods.

For the treatment of TNF- α , IL-1 β , IL-6, and IL-8 mediated diseases, cancer, and/or hyperglycemia, the compounds of the present invention may be administered orally, parentally, by inhalation spray, rectally, or topically in dosage unit formulations containing conventional pharmaceutically acceptable carriers, adjuvants, and vehicles. The term parenteral as used herein includes, subcutaneous, intravenous, intramuscular, intrasternal, infusion techniques or intraperitoneally.

The dosage regimen for treating a TNF- α , IL-1, IL-6, and IL-8 mediated diseases, cancer, and/or hyperglycemia with the compounds of this invention and/or compositions of this invention is based on a variety of factors, including the type of disease, the age, weight, sex, medical condition of the patient, the severity of the condition, the route of administration, and the particular compound employed. Thus, the dosage regimen may vary widely, but can be determined routinely using standard methods. Dosage levels of the order from about 0.01 mg to 30 mg per kilogram of body weight per day, preferably from about 0.1 mg to 10 mg/kg, more preferably from about 0.25 mg to 1 mg/kg are useful for all methods of use disclosed herein.

The pharmaceutically active compounds of this invention can be processed in accordance with conventional methods of pharmacy to produce medicinal agents for administration to patients, including humans and other mammals.

For oral administration, the pharmaceutical composition may be in the form of, for example, a capsule, a tablet, a suspension, or liquid. The pharmaceutical composition is preferably made in the form of a dosage unit containing a given amount of the active ingredient. For example, these may contain an

amount of active ingredient from about 1 to 2000 mg, preferably from about 1 to 500 mg, more preferably from about 5 to 150 mg. A suitable daily dose for a human or other mammal may vary widely depending on the condition of the patient and other factors, but, once again, can be determined using routine methods.

The active ingredient may also be administered by injection as a composition with suitable carriers including saline, dextrose, or water. The daily parenteral dosage regimen will be from about 0.1 to about 30 mg/kg of total body weight, preferably from about 0.1 to about 10 mg/kg, and more preferably from about 0.25 mg to 1 mg/kg.

Injectable preparations, such as sterile injectable aqueous or oleaginous suspensions, may be formulated according to the known art using suitable dispersing or wetting agents and suspending agents. The sterile injectable preparation may also be a sterile injectable solution or suspension in a non-toxic parenterally acceptable diluent or solvent, for example as a solution in 1,3-butanediol. Among the acceptable vehicles and solvents that may be employed are water, Ringer's solution, and isotonic sodium chloride solution. In addition, sterile, fixed oils are conventionally employed as a solvent or suspending medium. For this purpose any bland fixed oil may be employed, including synthetic mono- or diglycerides. In addition, fatty acids such as oleic acid find use in the preparation of injectables.

Suppositories for rectal administration of the drug can be prepared by mixing the drug with a suitable non-irritating excipient such as cocoa butter and polyethylene glycols that are solid at ordinary temperatures but liquid at the rectal temperature and will therefore melt in the rectum and release the drug.

A suitable topical dose of active ingredient of a compound of the invention is 0.1 mg to 150 mg

administered one to four, preferably one or two times daily. For topical administration, the active ingredient may comprise from 0.001% to 10% w/w, e.g., from 1% to 2% by weight of the formulation, although it
5 may comprise as much as 10% w/w, but preferably not more than 5% w/w, and more preferably from 0.1% to 1% of the formulation.

Formulations suitable for topical administration include liquid or semi-liquid preparations suitable for
10 penetration through the skin (e.g., liniments, lotions, ointments, creams, or pastes) and drops suitable for administration to the eye, ear, or nose.

For administration, the compounds of this invention are ordinarily combined with one or more adjuvants
15 appropriate for the indicated route of administration. The compounds may be admixed with lactose, sucrose, starch powder, cellulose esters of alkanolic acids, stearic acid, talc, magnesium stearate, magnesium oxide, sodium and calcium salts of phosphoric and sulphuric
20 acids, acacia, gelatin, sodium alginate, polyvinylpyrrolidone, and/or polyvinyl alcohol, and tableted or encapsulated for conventional administration. Alternatively, the compounds of this invention may be dissolved in saline, water,
25 polyethylene glycol, propylene glycol, ethanol, corn oil, peanut oil, cottonseed oil, sesame oil, tragacanth gum, and/or various buffers. Other adjuvants and modes of administration are well known in the pharmaceutical art. The carrier or diluent may include time delay
30 material, such as glyceryl monostearate or glyceryl distearate alone or with a wax, or other materials well known in the art.

The pharmaceutical compositions may be made up in a solid form (including granules, powders or
35 suppositories) or in a liquid form (e.g., solutions, suspensions, or emulsions). The pharmaceutical compositions may be subjected to conventional

pharmaceutical operations such as sterilization and/or may contain conventional adjuvants, such as preservatives, stabilizers, wetting agents, emulsifiers, buffers etc.

5 Solid dosage forms for oral administration may include capsules, tablets, pills, powders, and granules. In such solid dosage forms, the active compound may be admixed with at least one inert diluent such as sucrose, lactose, or starch. Such dosage forms may also
10 comprise, as in normal practice, additional substances other than inert diluents, e.g., lubricating agents such as magnesium stearate. In the case of capsules, tablets, and pills, the dosage forms may also comprise buffering agents. Tablets and pills can additionally be
15 prepared with enteric coatings.

 Liquid dosage forms for oral administration may include pharmaceutically acceptable emulsions, solutions, suspensions, syrups, and elixirs containing inert diluents commonly used in the art, such as water.
20 Such compositions may also comprise adjuvants, such as wetting, sweetening, flavoring, and perfuming agents.

 Compounds of the present invention can possess one or more asymmetric carbon atoms and are thus capable of existing in the form of optical isomers as well as in
25 the form of racemic or non-racemic mixtures thereof. The optical isomers can be obtained by resolution of the racemic mixtures according to conventional processes, e.g., by formation of diastereoisomeric salts, by treatment with an optically active acid or base.
30 Examples of appropriate acids are tartaric, diacetyltartaric, dibenzoyltartaric, ditoluoyltartaric, and camphorsulfonic acid and then separation of the mixture of diastereoisomers by crystallization followed by liberation of the optically active bases from these
35 salts. A different process for separation of optical isomers involves the use of a chiral chromatography column optimally chosen to maximize the separation of

the enantiomers. Still another available method involves synthesis of covalent diastereoisomeric molecules by reacting compounds of the invention with an optically pure acid in an activated form or an optically pure isocyanate. The synthesized diastereoisomers can be separated by conventional means such as chromatography, distillation, crystallization or sublimation, and then hydrolyzed to deliver the enantiomerically pure compound. The optically active compounds of the invention can likewise be obtained by using active starting materials. These isomers may be in the form of a free acid, a free base, an ester or a salt.

The compounds of the present invention can be used in the form of salts derived from inorganic or organic acids. The salts include, but are not limited to, the following: acetate, adipate, alginate, citrate, aspartate, benzoate, benzenesulfonate, bisulfate, butyrate, camphorate, camphorsulfonate, digluconate, cyclopentanepropionate, dodecylsulfate, ethanesulfonate, glucoheptanoate, glycerophosphate, hemisulfate, heptanoate, hexanoate, fumarate, hydrochloride, hydrobromide, hydroiodide, 2-hydroxy-ethanesulfonate, lactate, maleate, methanesulfonate, nicotinate, 2-naphthalenesulfonate, oxalate, palmoate, pectinate, persulfate, 2-phenylpropionate, picrate, pivalate, propionate, succinate, tartrate, thiocyanate, tosylate, mesylate, and undecanoate. Also, the basic nitrogen-containing groups can be quaternized with such agents as lower alkyl halides, such as methyl, ethyl, propyl, and butyl chloride, bromides and iodides; dialkyl sulfates like dimethyl, diethyl, dibutyl, and diamyl sulfates, long chain halides such as decyl, lauryl, myristyl and stearyl chlorides, bromides and iodides, aralkyl halides like benzyl and phenethyl bromides, and others. Water or oil-soluble or dispersible products are thereby obtained.

Examples of acids that may be employed to form pharmaceutically acceptable acid addition salts include such inorganic acids as hydrochloric acid, sulphuric acid and phosphoric acid and such organic acids as
5 oxalic acid, maleic acid, succinic acid and citric acid. Other examples include salts with alkali metals or alkaline earth metals, such as sodium, potassium, calcium or magnesium or with organic bases.

While the compounds of the invention can be
10 administered as the sole active pharmaceutical agent, they can also be used in combination with one or more compounds of the invention or other agents. When administered as a combination, the therapeutic agents can be formulated as separate compositions that are
15 given at the same time or different times, or the therapeutic agents can be given as a single composition.

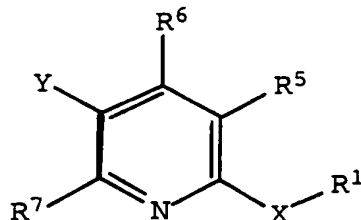
The foregoing is merely illustrative of the invention and is not intended to limit the invention to the disclosed compounds. Variations and changes which
20 are obvious to one skilled in the art are intended to be within the scope and nature of the invention which are defined in the appended claims.

From the foregoing description, one skilled in the art can easily ascertain the essential characteristics
25 of this invention, and without departing from the spirit and scope thereof, can make various changes and modifications of the invention to adapt it to various usages and conditions.

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WHAT IS CLAIMED IS:

1. A compound of the formula



5 or a pharmaceutically acceptable salt thereof, wherein

X is O, S, S(O), S(O)₂, or NR²;Y is -C(O)-NR³R⁴ or -NR⁴-C(O)-R³;

10

R¹ is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of alkyl, halo, haloalkyl, cyano, azido, nitro, amidino, R¹⁸-Z¹⁸- or R¹⁸-Z¹⁸-alkyl; provided that the total number of aryl, heteroaryl, cycloalkyl and heterocyclyl radicals in R¹ is 1-3; and provided when Y is -NR⁴-C(O)-R³ and X is O or S, R¹ is other than a 2-pyrimidinyl radical;

20 R² is a hydrogen or alkyl radical;

R³ is an aryl or heteroaryl radical which is optionally substituted by 1-5 radicals of alkyl, halo, haloalkyl, cyano, azido, nitro, amidino, R¹⁹-Z¹⁹- or R¹⁹-Z¹⁹-alkyl; provided that the total number of aryl and heteroaryl radicals in R³ is 1-3; and provided when Y is -C(O)-NR³R⁴, R³ is other than a phenyl or naphthyl having an amino, nitro, cyano, carboxy or alkoxycarbonyl

substituent bonded to the ring carbon atom adjacent to the ring carbon atom bonded to $-NR^4-$; and

R^4 is a hydrogen, alkyl, alkenyl, haloalkyl,
5 haloalkenyl, aryl, heteroaryl, arylalkyl,
heteroarylalkyl or $R^{20}-Z^{20}$ -alkyl radical;

wherein R^{18} , R^{19} and R^{20} are each independently a
hydrogen, alkyl, haloalkyl, aryl, heteroaryl, arylalkyl
10 or heteroarylalkyl radical; wherein the aryl and
heteroaryl radicals of R^4 , R^{18} , R^{19} and R^{20} are optionally
substituted by 1-3 radicals of hydroxy, alkoxy,
alkylthiol, amino, alkylamino, dialkylamino,
alkanoylamino, alkylsulfonylamino, alkylsulfinyl,
15 alkylsulfonyl, alkoxycarbonylamino, alkoxycarbonyl,
cyano, halo, azido, alkyl, haloalkyl or haloalkoxy; and

Z^{18} , Z^{19} and Z^{20} are each independently $-O-$, $-S-$, $-S(O)-$,
 $-S(O)_2-$, $-CO_2-$, $-C(O)-$, $-NR^{21}-$, $-NR^{21}-C(O)-$, $-C(O)-NR^{21}-$,
20 $NR^{21}-S(O)_2-$ or $-S(O)_2-NR^{21}-$; wherein each R^{21} is
independently a hydrogen or alkyl radical;

R^5 and R^6 are each independently a hydrogen, alkyl, halo,
haloalkyl, haloalkoxy, aminoalkyl, alkylaminoalkyl,
25 dialkylaminoalkyl, amino, alkylamino, dialkylamino,
alkanoylamino, alkylsulfonylamino, aminosulfonyl,
alkylaminosulfonyl, dialkylaminosulfonyl, hydroxy,
hydroxyalkyl, thiol, alkylthiol, alkylsulfinyl,
alkylsulfonyl, alkoxy, alkoxyalkyl, cyano, azido, nitro,
30 carboxy, alkoxycarbonyl, aminocarbonyl,
alkylaminocarbonyl or dialkylaminocarbonyl radical; and

R^7 is a hydrogen, alkyl, halo, haloalkyl, haloalkoxy,
aminoalkyl, alkylaminoalkyl, dialkylaminoalkyl,

aminosulfonyl, alkylaminosulfonyl, dialkylaminosulfonyl, hydroxy, hydroxyalkyl, thiol, alkylthiol, alkylsulfinyl, alkylsulfonyl, alkoxy, alkoxyalkyl, cyano, azido, nitro, carboxy, alkoxycarbonyl, aminocarbonyl,
5 alkylaminocarbonyl or dialkylaminocarbonyl radical.

2. The compound of Claim 1 or a pharmaceutically acceptable salt thereof, wherein
10 R^1 is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of C_1 - C_6 alkyl, halo, C_1 - C_6 haloalkyl of 1-3 halo radicals, cyano, azido, nitro, amidino, R^{18} - Z^{18} - or R^{18} -
15 Z^{18} - C_1 - C_6 alkyl; provided that the total number of aryl, heteroaryl, cycloalkyl and heterocyclyl radicals in R^1 is 1-3; and provided when Y is $-NR^4-C(O)-R^3$ and X is O or S, R^1 is other than a 2-pyrimidinyl radical;

20 R^2 is a hydrogen or C_1 - C_4 alkyl radical;

R^3 is an aryl or heteroaryl radical which is optionally substituted by 1-5 radicals of C_1 - C_6 alkyl, halo, C_1 - C_6 haloalkyl of 1-3 halo radicals, cyano, azido, nitro,
25 amidino, R^{19} - Z^{19} - or R^{19} - Z^{19} - C_1 - C_6 alkyl; provided that the total number of aryl and heteroaryl radicals in R^3 is 1-3; and provided when Y is $-C(O)-NR^3R^4$, R^3 is other than a phenyl or naphthyl having an amino, nitro, cyano, carboxy or alkoxycarbonyl substituent bonded to the ring
30 carbon atom adjacent to the ring carbon atom bonded to $-NR^4$; and

R^4 is a radical of hydrogen, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_1 - C_6 haloalkyl of 1-3 halo radicals, C_2 - C_6 haloalkenyl of 1-3 halo radicals, aryl, heteroaryl, aryl- C_1 - C_4 alkyl, heteroaryl- C_1 - C_4 alkyl or R^{20} - Z^{20} - C_1 - C_6 alkyl radical; and

wherein R^{18} , R^{19} and R^{20} are each independently a hydrogen, C_1 - C_4 alkyl, C_1 - C_4 haloalkyl of 1-3 halo radicals, aryl, heteroaryl, aryl- C_1 - C_4 alkyl or heteroaryl- C_1 - C_4 alkyl radical; wherein the aryl and heteroaryl radicals of R^4 , R^{18} , R^{19} and R^{20} are optionally substituted by 1-3 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, C_1 - C_5 alkanoylamino, C_1 - C_4 alkylsulfonylamino, C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkylsulfonyl, (C_1 - C_4 alkoxy)carbonylamino, (C_1 - C_4 alkoxy)carbonyl, cyano, halo, azido, C_1 - C_4 alkyl, C_1 - C_4 haloalkyl of 1-3 halo radicals or C_1 - C_4 haloalkoxy of 1-3 halo radicals; and

each R^{21} is independently a hydrogen or C_1 - C_4 alkyl radical;

R^5 and R^6 are each independently a hydrogen, C_1 - C_4 alkyl, halo, C_1 - C_4 haloalkyl of 1-3 halo radicals, C_1 - C_4 haloalkoxy of 1-3 halo radicals, C_1 - C_4 aminoalkyl, (C_1 - C_4 alkyl)amino- C_1 - C_4 alkyl, di(C_1 - C_4 alkyl)amino- C_1 - C_4 alkyl, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, C_1 - C_5 alkanoylamino, C_1 - C_4 alkylsulfonylamino, aminosulfonyl, C_1 - C_4 alkylaminosulfonyl, di(C_1 - C_4 alkyl)aminosulfonyl, hydroxy, C_1 - C_4 hydroxyalkyl, thiol, C_1 - C_4 alkylthiol, C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkylsulfonyl, C_1 - C_4 alkoxy, (C_1 - C_4 alkoxy) C_1 - C_4 alkyl,

cyano, azido, nitro, carboxy, (C₁-C₄ alkoxy)carbonyl, aminocarbonyl, (C₁-C₄ alkyl)aminocarbonyl or di(C₁-C₄ alkyl)aminocarbonyl radical; and

- 5 R⁷ is a hydrogen, C₁-C₄ alkyl, halo, C₁-C₄ haloalkyl of 1-3 halo radicals, C₁-C₄ haloalkoxy of 1-3 halo radicals, C₁-C₄ aminoalkyl, (C₁-C₄ alkyl)amino-C₁-C₄ alkyl, di(C₁-C₄ alkyl)amino-C₁-C₄ alkyl, aminosulfonyl, C₁-C₄ alkylaminosulfonyl, di(C₁-C₄ alkyl)aminosulfonyl,
10 hydroxy, C₁-C₄ hydroxyalkyl, thiol, C₁-C₄ alkylthiol, C₁-C₄ alkylsulfinyl, C₁-C₄ alkylsulfonyl, C₁-C₄ alkoxy, (C₁-C₄ alkoxy)C₁-C₄ alkyl, cyano, azido, nitro, carboxy, (C₁-C₄ alkoxy)carbonyl, aminocarbonyl, (C₁-C₄ alkyl)aminocarbonyl or di(C₁-C₄ alkyl)aminocarbonyl
15 radical; and

wherein cycloalkyl is a monocyclic, bicyclic or tricyclic carbocyclic alkyl radical of 5-12 ring members, which is optionally partially unsaturated,
20 benzo fused or heterocyclo fused; aryl is a phenyl or biphenyl radical which is optionally benzo fused or heterocyclo fused; heterocyclyl is a radical of a monocyclic or bicyclic saturated heterocyclic ring system having 5-8 ring members per ring, wherein 1-3
25 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally partially unsaturated or benzo-fused and optionally substituted by 1-2 oxo or thioxo radicals; and heteroaryl is a monocyclic or bicyclic aromatic heterocyclic ring system having 5-6 ring
30 members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C₃-C₄-carbocyclic-fused.

- 35 3. The compound of Claim 2 or a pharmaceutically acceptable salt thereof, wherein Y is -NR⁴-C(O)-R³.

4. The compound of Claim 3 or a pharmaceutically acceptable salt thereof, wherein

5

X is O or NR^2 ;

R^1 is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of $\text{C}_1\text{-C}_4$ alkyl, halo, $\text{C}_1\text{-C}_4$ haloalkyl of 1-3 halo radicals, cyano, azido, nitro, amidino, $\text{R}^{18}\text{-Z}^{18}\text{-}$ or $\text{R}^{18}\text{-Z}^{18}\text{-C}_1\text{-C}_4$ alkyl; provided that the total number of aryl, heteroaryl, cycloalkyl and heterocyclyl radicals in R^1 is 1-2;

15

wherein each R^{18} is independently a hydrogen, $\text{C}_1\text{-C}_4$ alkyl, trifluoromethyl, aryl, heteroaryl, aryl- $\text{C}_1\text{-C}_2$ alkyl or heteroaryl- $\text{C}_1\text{-C}_2$ alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, $\text{C}_1\text{-C}_4$ alkoxy, $\text{C}_1\text{-C}_4$ alkylthiol, amino, $\text{C}_1\text{-C}_4$ alkylamino, di($\text{C}_1\text{-C}_4$ alkyl)amino, acetylamino, cyano, halo, azido, $\text{C}_1\text{-C}_4$ alkyl, trifluoromethyl or trifluoromethoxy; and

25 R^2 is a hydrogen or $\text{C}_1\text{-C}_2$ alkyl radical;

R^3 is an aryl or heteroaryl radical which is optionally substituted by 1-5 radicals of $\text{C}_1\text{-C}_6$ alkyl, halo, $\text{C}_1\text{-C}_4$ haloalkyl of 1-3 halo radicals, cyano, azido, nitro, amidino, $\text{R}^{19}\text{-Z}^{19}\text{-}$ or $\text{R}^{19}\text{-Z}^{19}\text{-C}_1\text{-C}_4$ alkyl; provided that the total number of aryl and heteroaryl radicals in R^3 is 1-2; and

30

wherein each R^{19} is independently a hydrogen, C_1 - C_4 alkyl, trifluoromethyl, aryl, heteroaryl, aryl- C_1 - C_4 alkyl or heteroaryl- C_1 - C_4 alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted
5 by 1-2 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, acetylamino, cyano, halo, C_1 - C_4 alkyl, trifluoromethyl or trifluoromethoxy; and

10 R^4 is a radical of hydrogen, C_1 - C_6 alkyl, aryl, heteroaryl, aryl- C_1 - C_4 alkyl, heteroaryl- C_1 - C_4 alkyl or R^{20} - Z^{20} - C_2 - C_4 alkyl radical; and

wherein R^{20} is a hydrogen, C_1 - C_4 alkyl, aryl, heteroaryl,
15 aryl- C_1 - C_2 alkyl or heteroaryl- C_1 - C_2 alkyl radical;
wherein the aryl and heteroaryl radicals of R^4 and R^{20} are optionally substituted by 1-2 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, acetylamino, halo, C_1 - C_4 alkyl,
20 trifluoromethyl or trifluoromethoxy; and

Z^{20} is -O- or - NR^{21} -; wherein each R^{21} is independently a hydrogen or methyl radical;

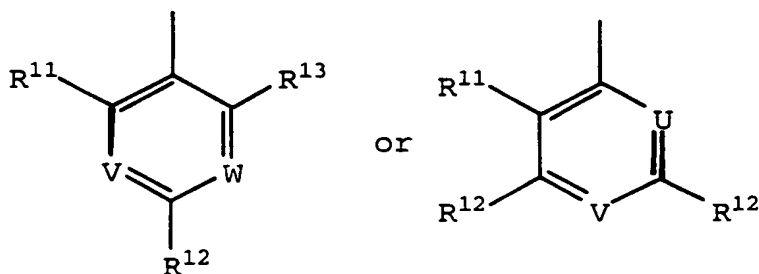
25 R^5 and R^6 are each independently a hydrogen, C_1 - C_4 alkyl, halo, trifluoromethyl, trifluoromethoxy, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, C_1 - C_5 alkanoylamino, hydroxy, C_1 - C_4 hydroxyalkyl, C_1 - C_4 alkoxy, cyano, azido, nitro, carboxy, (C_1 - C_4 alkoxy)carbonyl, aminocarbonyl,
30 (C_1 - C_4 alkyl)aminocarbonyl or di(C_1 - C_4 alkyl)aminocarbonyl radical; and

R^7 is a hydrogen, C_1 - C_4 alkyl, halo, trifluoromethyl, trifluoromethoxy, hydroxy, C_1 - C_4 hydroxyalkyl, C_1 - C_4 alkoxy, carboxy, (C_1 - C_4 alkoxy)carbonyl, aminocarbonyl, (C_1 - C_4 alkyl)aminocarbonyl or di(C_1 - C_4 alkyl)aminocarbonyl radical; and

wherein cycloalkyl is a monocyclic or bicyclic carbocyclic alkyl radical of 5-12 ring members, which is optionally partially unsaturated, benzo fused or heterocyclo fused; aryl is a phenyl or biphenyl radical which is optionally benzo fused or heterocyclo fused; heterocyclyl is a radical of a monocyclic or bicyclic saturated heterocyclic ring system having 5-8 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally partially unsaturated or benzo-fused and optionally substituted by 1-2 oxo or thioxo radicals; and heteroaryl is a monocyclic aromatic heterocyclic ring system having 5-6 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C_3 - C_4 -carbocyclic-fused.

5. The compound of Claim 4 or a pharmaceutically acceptable salt thereof, wherein

R^3 is a radical of the formula



wherein

U is C-R¹³ or N;

V and W are each independently C-R¹² or N;

5

R¹¹ and R¹³ are each independently a radical of hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl, cyano, azido, nitro, amidino or R¹⁹-Z¹⁹-; and each R¹² is independently a radical of hydrogen, C₁-C₆ alkyl, halo, C₁-C₄ haloalkyl
10 of 1-3 halo radicals, R³¹-Z³¹- or R³¹-Z³¹-C₁-C₄ alkyl; provided that the combined total number of aryl and heteroaryl radicals in R¹¹, R¹² and R¹³ is 0-1;

wherein each R¹⁹ is independently a hydrogen, C₁-C₄
15 alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, halo,
20 methyl, ethyl, trifluoromethyl or trifluoromethoxy; and

each Z¹⁹ is independently -O-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-;

25 wherein each R³¹ is independently a hydrogen, C₁-C₄ alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino,
30 methylamino, dimethylamino, acetylamino, cyano, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy; and

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each Z^{31} is independently -O-, $-NR^{21}-$, $-NR^{21}-C(O)-$, $-C(O)-NR^{21}-$, $-NR^{21}-S(O)_2-$ or $-S(O)_2-NR^{21}-$;

wherein R^4 is a radical of hydrogen, C_1-C_6 alkyl, aryl,
5 heteroaryl, aryl- C_1-C_4 alkyl, heteroaryl- C_1-C_4 alkyl or
 $R^{20}-Z^{20}-C_2-C_4$ alkyl radical;

wherein R^{20} is a hydrogen, C_1-C_4 alkyl, aryl, heteroaryl,
aryl- C_1-C_2 alkyl or heteroaryl- C_1-C_2 alkyl radical;
10 wherein the aryl and heteroaryl radicals of R^4 and R^{20}
are optionally substituted by 1-2 radicals of hydroxy,
methoxy, ethoxy, methylthiol, ethylthiol, amino,
methylamino, dimethylamino, ethylamino, diethylamino,
acetylamino, halo, methyl, ethyl, trifluoromethyl or
15 trifluoromethoxy; and

R^5 and R^6 are each independently a hydrogen, methyl,
ethyl, halo, trifluoromethyl, trifluoromethoxy, amino,
 C_1-C_2 alkylamino, di(C_1-C_2 alkyl)amino, hydroxy, methoxy
20 or ethoxy radical; and

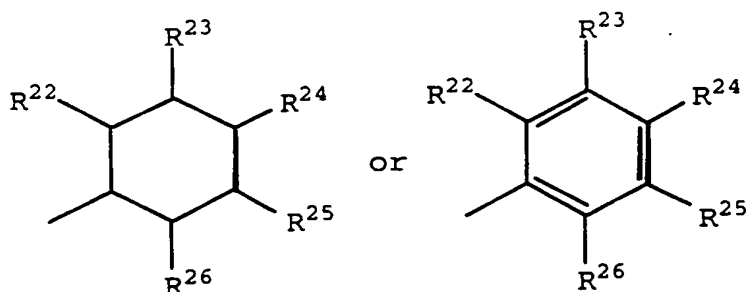
R^7 is a hydrogen, methyl, ethyl, halo, trifluoromethyl,
trifluoromethoxy, hydroxy, methoxy or ethoxy radical;
and

25 wherein cycloalkyl is a monocyclic or bicyclic
carbocyclic alkyl radical of 5-10 ring members, which is
optionally partially unsaturated with one double bond,
benzo fused or heterocyclo fused; aryl is a phenyl or
30 biphenyl radical which is optionally benzo fused or
heterocyclo fused; heterocyclyl is a radical of a
monocyclic or bicyclic saturated heterocyclic ring
system having 5-8 ring members per ring, wherein 1-3
ring members are oxygen, sulfur or nitrogen heteroatoms,

which is optionally partially unsaturated or benzo-fused and optionally substituted by 1-2 oxo or thioxo radicals; and heteroaryl is a monocyclic aromatic heterocyclic ring system having 5-6 ring members per
 5 ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C₃-C₄-carbocyclic-fused.

10 6. The compound of Claim 5 or a pharmaceutically acceptable salt thereof, wherein

R¹ is a radical of the formula



15 wherein

R²², R²³, R²⁴, R²⁵ and R²⁶ are each independently a radical of hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl, cyano, azido, nitro, amidino, R¹⁸-Z¹⁸- or R¹⁸-Z¹⁸-C₁-C₄ alkyl;
 20 provided at least one of R²¹, R²², R²³, R²⁴ and R²⁵ is hydrogen; and provided that the combined total number of aryl and heteroaryl radicals in R²², R²³, R²⁴, R²⁵ and R²⁶ is 0-1;

25 wherein each Z¹⁸ is independently -O-, -S-, -S(O)₂-, -CO₂-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; and

wherein aryl is a phenyl or biphenyl radical which is optionally benzo fused or heterocyclo fused; and heteroaryl is a monocyclic aromatic heterocyclic ring system having 5-6 ring members per ring, wherein 1-3
5 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C₃-C₄-carbocyclic-fused.

10 7. The compound of Claim 6 or a pharmaceutically acceptable salt thereof, wherein

X is NR²;

15 R² is a hydrogen or methyl radical;

R⁴ is a radical of hydrogen, methyl or ethyl radical;
and

20 R⁵, R⁶ and R⁷ are each independently a hydrogen radical;
and

wherein aryl is a phenyl, biphenyl or naphthyl radical;
and heteroaryl is a monocyclic aromatic heterocyclic
25 ring system having 5-6 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms.

30 8. The compound of Claim 7 or a pharmaceutically acceptable salt thereof, wherein

R¹¹ and R¹³ are each independently a radical of hydrogen, methyl, ethyl, fluoro, chloro, trifluoromethyl, cyano,
35 azido, nitro, amidino, R¹⁹-O-, R¹⁹-S(O)₂-, R¹⁹-O-C(O)-,

- $R^{19}-C(O)-$, $R^{19}-NR^{21}-C(O)-$ or $R^{19}-NR^{21}-S(O)_2-$; and each R^{12} is independently a radical of hydrogen, methyl, ethyl, fluoro, chloro, trifluoromethyl, trifluoromethoxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, aminocarbonyl, methylaminocarbonyl, dimethylaminocarbonyl, aminomethyl, (methylamino)methyl or (dimethylamino)methyl; provided that the combined total number of aryl and heteroaryl radicals in R^{11} , R^{12} and R^{13} is 0-1; and
- wherein each R^{19} is independently a hydrogen, methyl, ethyl, trifluoromethyl, phenyl, heteroaryl, phenylmethyl or heteroaryl-methyl radical; wherein the phenyl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, fluoro, chloro, methyl, ethyl, trifluoromethyl or trifluoromethoxy.
9. The compound of Claim 2 or a pharmaceutically acceptable salt thereof, wherein Y is $-C(O)-NR^3R^4$.
10. The compound of Claim 9 or a pharmaceutically acceptable salt thereof, wherein
- X is O or NR^2 ;
- R^1 is a cycloalkyl, aryl, heterocyclyl or heteroaryl radical which is optionally substituted by 1-4 radicals of C_1-C_4 alkyl, halo, C_1-C_4 haloalkyl of 1-3 halo radicals, cyano, azido, nitro, amidino, $R^{18}-Z^{18}-$ or $R^{18}-Z^{18}-C_1-C_4$ alkyl; provided that the total number of aryl,

heteroaryl, cycloalkyl and heterocyclyl radicals in R^1 is 1-2;

wherein each R^{18} is independently a hydrogen, C_1 - C_4 alkyl, trifluoromethyl, aryl, heteroaryl, aryl- C_1 - C_2 alkyl or heteroaryl- C_1 - C_2 alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, C_1 - C_4 alkoxy, C_1 - C_4 alkylthiol, amino, C_1 - C_4 alkylamino, di(C_1 - C_4 alkyl)amino, acetylamino, cyano, halo, azido, C_1 - C_4 alkyl, trifluoromethyl or trifluoromethoxy; and

each Z^{18} is independently -O-, -S-, -S(O)-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; wherein each R^{21} is independently a hydrogen or C_1 - C_4 alkyl radical;

R^2 is a hydrogen or C_1 - C_2 alkyl radical;

R^3 is an aryl or heteroaryl radical which is optionally substituted by 1-5 radicals of C_1 - C_6 alkyl, halo, C_1 - C_4 haloalkyl of 1-3 halo radicals, cyano, azido, nitro, amidino, R^{19} -Z¹⁹- or R^{19} -Z¹⁹- C_1 - C_4 alkyl; provided that the total number of aryl and heteroaryl radicals in R^3 is 1-2; and provided R^3 is other than a phenyl or naphthyl having an amino, nitro, cyano, carboxy or alkoxycarbonyl substituent bonded to the ring carbon atom adjacent to the ring carbon atom bonded to -NR⁴-; and

wherein each R^{19} is independently a hydrogen, C_1 - C_4 alkyl, trifluoromethyl, aryl, heteroaryl, aryl- C_1 - C_4 alkyl or heteroaryl- C_1 - C_4 alkyl radical; wherein the

aryl and heteroaryl radicals are optionally substituted
by 1-2 radicals of hydroxy, C₁-C₄ alkoxy, C₁-C₄
alkylthiol, amino, C₁-C₄ alkylamino, di(C₁-C₄
alkyl)amino, acetylamino, cyano, halo, C₁-C₄ alkyl,
5 trifluoromethyl or trifluoromethoxy; and

R⁴ is a radical of hydrogen, C₁-C₆ alkyl, aryl,
heteroaryl, aryl-C₁-C₄ alkyl, heteroaryl-C₁-C₄ alkyl or
R²⁰-Z²⁰-C₂-C₄ alkyl radical; and

10 wherein R²⁰ is a hydrogen, C₁-C₄ alkyl, aryl, heteroaryl,
aryl-C₁-C₂ alkyl or heteroaryl-C₁-C₂ alkyl radical;
wherein the aryl and heteroaryl radicals of R⁴ and R²⁰
are optionally substituted by 1-2 radicals of hydroxy,
15 C₁-C₄ alkoxy, C₁-C₄ alkylthiol, amino, C₁-C₄ alkylamino,
di(C₁-C₄ alkyl)amino, acetylamino, halo, C₁-C₄ alkyl,
trifluoromethyl or trifluoromethoxy; and

Z²⁰ is -O- or -NR²¹-; wherein each R²¹ is independently a
20 hydrogen or methyl radical;

R⁵ and R⁶ are each independently a hydrogen, C₁-C₄ alkyl,
halo, trifluoromethyl, trifluoromethoxy, amino, C₁-C₄
alkylamino, di(C₁-C₄ alkyl)amino, C₁-C₅ alkanoylamino,
25 hydroxy, C₁-C₄ hydroxyalkyl, C₁-C₄ alkoxy, cyano, azido,
nitro, carboxy, (C₁-C₄ alkoxy)carbonyl, aminocarbonyl,
(C₁-C₄ alkyl)aminocarbonyl or di(C₁-C₄
alkyl)aminocarbonyl radical; and

30 R⁷ is a hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl,
trifluoromethoxy, hydroxy, C₁-C₄ hydroxyalkyl, C₁-C₄
alkoxy, carboxy, (C₁-C₄ alkoxy)carbonyl, aminocarbonyl,

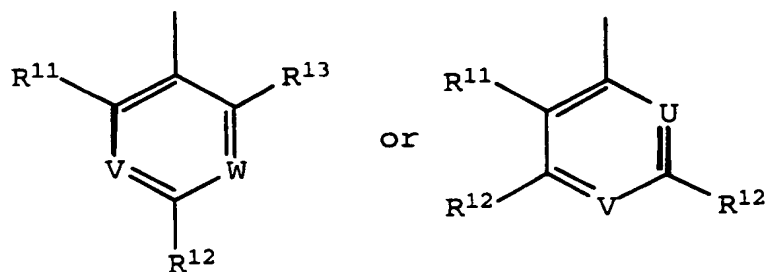
(C₁-C₄ alkyl)aminocarbonyl or di(C₁-C₄ alkyl)aminocarbonyl radical; and

wherein cycloalkyl is a monocyclic or bicyclic
 5 carbocyclic alkyl radical of 5-12 ring members, which is optionally partially unsaturated, benzo fused or heterocyclo fused; aryl is a phenyl or biphenyl radical which is optionally benzo fused or heterocyclo fused; heterocyclyl is a radical of a monocyclic or bicyclic
 10 saturated heterocyclic ring system having 5-8 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally partially unsaturated or benzo-fused and optionally substituted by 1-2 oxo or thioxo radicals; and
 15 heteroaryl is a monocyclic aromatic heterocyclic ring system having 5-6 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C₃-C₄-carbocyclic-fused.

20

11. The compound of Claim 10 or a pharmaceutically acceptable salt thereof, wherein

25 R³ is a radical of the formula



wherein

U is C-R¹³ or N;

30

V and W are each independently C-R¹² or N;

R¹¹ and R¹³ are each independently a radical of hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl, cyano, azido, nitro, amidino or R¹⁹-Z¹⁹-; and each R¹² is independently a radical of hydrogen, C₁-C₆ alkyl, halo, C₁-C₄ haloalkyl of 1-3 halo radicals, R³¹-Z³¹- or R³¹-Z³¹-C₁-C₄ alkyl; provided that the combined total number of aryl and heteroaryl radicals in R¹¹, R¹² and R¹³ is 0-1; provided when U is C-R¹³ and V and W are each C-R¹², R¹¹ and R¹³ are each other than a nitro, cyano, carboxy or alkoxycarbonyl radical;

wherein each R¹⁹ is independently a hydrogen, C₁-C₄ alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy; and

each Z¹⁹ is independently -O-, -S(O)₂-, -CO₂-, -C(O)-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; wherein each R²¹ is independently a hydrogen or methyl radical;

wherein each R³¹ is independently a hydrogen, C₁-C₄ alkyl, trifluoromethyl, aryl, heteroaryl, aryl-C₁-C₄ alkyl or heteroaryl-C₁-C₄ alkyl radical; wherein the aryl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, halo, methyl, ethyl, trifluoromethyl or trifluoromethoxy; and

each Z^{31} is independently -O-, $-NR^{21}-$, $-NR^{21}-C(O)-$, $-C(O)-NR^{21}-$, $-NR^{21}-S(O)_2-$ or $-S(O)_2-NR^{21}-$;

- 5 R^4 is a radical of hydrogen, C_1-C_6 alkyl, aryl, heteroaryl, aryl- C_1-C_4 alkyl, heteroaryl- C_1-C_4 alkyl or $R^{20}-Z^{20}-C_2-C_4$ alkyl radical; and

- wherein R^{20} is a hydrogen, C_1-C_4 alkyl, aryl, heteroaryl,
10 aryl- C_1-C_2 alkyl or heteroaryl- C_1-C_2 alkyl radical;
wherein the aryl and heteroaryl radicals of R^4 and R^{20}
are optionally substituted by 1-2 radicals of hydroxy,
methoxy, ethoxy, methylthiol, ethylthiol, amino,
methylamino, dimethylamino, ethylamino, diethylamino,
15 acetylamino, halo, methyl, ethyl, trifluoromethyl or
trifluoromethoxy; and

- R^5 and R^6 are each independently a hydrogen, methyl,
ethyl, halo, trifluoromethyl, trifluoromethoxy, amino,
20 C_1-C_2 alkylamino, di(C_1-C_2 alkyl)amino, hydroxy, methoxy
or ethoxy radical; and

- R^7 is a hydrogen, methyl, ethyl, halo, trifluoromethyl,
trifluoromethoxy, hydroxy, methoxy or ethoxy radical;
25 and

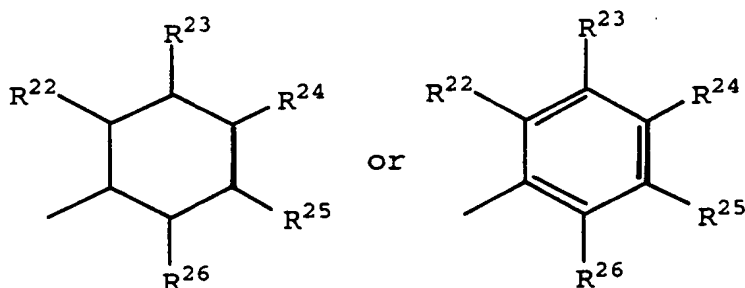
- wherein cycloalkyl is a monocyclic or bicyclic
carbocyclic alkyl radical of 5-10 ring members, which is
optionally partially unsaturated with one double bond,
30 benzo fused or heterocyclo fused; aryl is a phenyl or
biphenyl radical which is optionally benzo fused or
heterocyclo fused; heterocyclyl is a radical of a
monocyclic or bicyclic saturated heterocyclic ring
system having 5-8 ring members per ring, wherein 1-3

ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally partially unsaturated or benzo-fused and optionally substituted by 1-2 oxo or thioxo radicals; and heteroaryl is a monocyclic aromatic heterocyclic ring system having 5-6 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C₃-C₄-carbocyclic-fused.

10

12. The compound of Claim 11 or a pharmaceutically acceptable salt thereof, wherein

R¹ is a radical of the formula



15

wherein

R²², R²³, R²⁴, R²⁵ and R²⁶ are each independently a radical of hydrogen, C₁-C₄ alkyl, halo, trifluoromethyl, cyano, azido, nitro, amidino, R¹⁸-Z¹⁸- or R¹⁸-Z¹⁸-C₁-C₄ alkyl; provided at least one of R²¹, R²², R²³, R²⁴ and R²⁵ is hydrogen; and provided that the combined total number of aryl and heteroaryl radicals in R²², R²³, R²⁴, R²⁵ and R²⁶ is 0-1;

25

each Z¹⁸ is independently -O-, -S-, -S(O)₂-, -CO₂-, -NR²¹-, -NR²¹-C(O)-, -C(O)-NR²¹-, -NR²¹-S(O)₂- or -S(O)₂-NR²¹-; and

wherein aryl is a phenyl or biphenyl radical which is optionally benzo fused or heterocyclo fused; and heteroaryl is a monocyclic aromatic heterocyclic ring system having 5-6 ring members per ring, wherein 1-3
5 ring members are oxygen, sulfur or nitrogen heteroatoms, which is optionally benzo-fused or saturated C₃-C₄-carbocyclic-fused.

10 13. The compound of Claim 12 or a pharmaceutically acceptable salt thereof, wherein

X is NR²;

15 R² is a hydrogen or methyl radical;

R⁴ is a radical of hydrogen, methyl or ethyl radical;
and

20 R⁵, R⁶ and R⁷ are each independently a hydrogen radical;
and

wherein aryl is a phenyl, biphenyl or naphthyl radical;
and heteroaryl is a monocyclic aromatic heterocyclic
25 ring system having 5-6 ring members per ring, wherein 1-3 ring members are oxygen, sulfur or nitrogen heteroatoms.

30 14. The compound of Claim 13 or a pharmaceutically acceptable salt thereof, wherein

R¹¹ and R¹³ are each independently a radical of hydrogen, methyl, ethyl, fluoro, chloro, trifluoromethyl, cyano,
35 azido, nitro, amidino, R¹⁹-O-, R¹⁹-S(O)₂-, R¹⁹-O-C(O)-,

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$R^{19}-C(O)-$, $R^{19}-NR^{21}-C(O)-$ or $R^{19}-NR^{21}-S(O)_2-$; provided when U is $C-R^{13}$ and V and W are each $C-R^{12}$, R^{11} and R^{13} are each other than a nitro, cyano, carboxy or alkoxycarbonyl radical; and provided that the combined total number of
5 aryl and heteroaryl radicals in R^{11} and R^{13} is 0-1;

each R^{12} is independently a radical of hydrogen, methyl, ethyl, fluoro, chloro, trifluoromethyl, trifluoromethoxy, methoxy, ethoxy, amino, methylamino,
10 dimethylamino, acetylamino, aminocarbonyl, methylaminocarbonyl, dimethylaminocarbonyl, aminomethyl, (methylamino)methyl or (dimethylamino)methyl; and

wherein each R^{19} is independently a hydrogen, methyl, ethyl, trifluoromethyl, phenyl, heteroaryl, phenylmethyl
15 or heteroaryl-methyl radical; wherein the phenyl and heteroaryl radicals are optionally substituted by 1-2 radicals of hydroxy, methoxy, ethoxy, amino, methylamino, dimethylamino, acetylamino, cyano, fluoro, chloro, methyl, ethyl, trifluoromethyl or
20 trifluoromethoxy.

15. The compound of Claim 1 which is:
25

2-cyclohexyloxy-5-(2-chlorophenylcarbonylamino)pyridine;
2-cyclohexyloxy-5-(2-methylphenylcarbonylamino)pyridine;
2-cyclohexyloxy-5-(2,6-dichlorophenylcarbonylamino)
pyridine;
30 2-cyclohexyloxy-5-(2,6-dimethylphenylcarbonylamino)
pyridine;
2-(2,4-dimethylphenoxy)-5-(2-chlorophenylcarbonylamino)
pyridine;
2-(2,4-dimethylphenoxy)-5-(2,6-dichlorophenylcarbonyl
35 amino)pyridine;

- 2-(2,4-dimethylphenoxy)-5-(2-methylphenylcarbonylamino)
pyridine;
- 2-(2,6-dimethyl-4-chlorophenoxy)-5-(2,6-dimethylphenyl
carbonylamino) pyridine;
- 5 2-(2-methyl-4-fluorophenoxy)-5-(2-methylphenylcarbonyl
amino)pyridine;
- 2-(2-methyl-4-chlorophenoxy)-5-(2-chlorophenylcarbonyl
amino)pyridine;
- 2-(2-methyl-4-chlorophenoxy)-5-(2-methylphenylcarbonyl
10 amino)pyridine;
- 2-(2-methylphenoxy)-5-(2-chlorophenylcarbonylamino)
pyridine;
- 2-(2-methylphenoxy)-5-(2,6-dichlorophenyl
carbonylamino)pyridine;
- 15 2-(2-methylphenoxy)-5-(2-methylphenylcarbonyl
amino)pyridine;
- 2-(2-methyl-4-chlorophenoxy)-5-(2,6-dichlorophenyl
carbonylamino)pyridine;
- 2-(2-methyl-4-chlorophenoxy)-5-(2,6-dimethylphenyl
20 carbonylamino)pyridine;
- 2-(4-chlorophenoxy)-5-(2,6-dimethylphenylcarbonylamino)
pyridine;
- 2-(2-methyl-4-fluorophenoxy)-5-(2,6-dichlorophenyl
carbonylamino)pyridine;
- 25 2-(2-methyl-4-fluorophenoxy)-5-(2,6-dimethylphenyl
carbonylamino)pyridine;
- 2-(2-methylphenoxy)-5-(2,6-dimethylphenyl
carbonylamino)pyridine;
- 2-(2-methyl-4-fluorophenoxy)-5-(2-fluorophenylcarbonyl
30 amino)pyridine;
- 2-(2,4-dimethylphenoxy)-5-(2,6-dimethylphenylcarbonyl
amino)pyridine;
- 2-(1-naphthyloxy)-5-(2-methylphenylcarbonylamino)
pyridine;
- 35 2-(1-naphthyloxy)-5-(2,6-dichlorophenylcarbonylamino)
pyridine;

- 2-(1-naphthyloxy)-5-(2,6-dimethylphenylcarbonylamino)pyridine;
- 2-(2-methyl-3-pyridyloxy)-5-(2,6-dichlorophenylcarbonylamino)pyridine;
- 5 2-(2-methyl-4-chlorophenoxy)-5-((3,5-dimethyl-4-isoxazolyl)carbonylamino)pyridine;
- 2-(2-methyl-4-chlorophenylthiol)-5-(2-methylphenylcarbonyl amino)pyridine;
- 2-(2-methyl-4-chlorophenylthiol)-5-(2,6-
- 10 dimethylphenylcarbonyl amino)pyridine;
- 2-cyclohexylamino-5-(2,6-dichlorophenylcarbonylamino)pyridine;
- 2-cyclohexylamino-5-(2,6-dimethylphenylcarbonylamino)pyridine;
- 15 2-(2-methylcyclohexylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
- 2-(2-methylcyclohexylamino)-5-(2-methylphenylcarbonyl amino)pyridine;
- 2-(2,4-dimethylphenylamino)-5-(2-fluorophenylcarbonyl
- 20 amino)pyridine;
- 2-(2,4-dimethylphenylamino)-5-(2-chlorophenylcarbonyl amino)pyridine;
- 2-(2,4-dimethylphenylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
- 25 2-(2-methyl-4-chlorophenylamino)-5-(2,6-dichlorophenylcarbonylamino)pyridine;
- 2-(2,4-dimethylphenylamino)-5-(2-methylphenylcarbonyl amino)pyridine;
- 2-(2-methylphenylamino)-5-(2-methylphenylcarbonyl
- 30 amino)pyridine;
- 2-(2-methylphenylamino)-5-(2,6-dichlorophenylcarbonyl amino)pyridine;
- 2-(2-methylphenylamino)-5-(2,6-dimethylphenylcarbonyl amino)pyridine;
- 35 2-(2,4-dimethylphenylamino)-5-(2,6-dimethylphenylcarbonyl amino)pyridine;

2-(2-methyl-4-chlorophenylamino)-5-(2-methylphenyl
carbonylamino)pyridine;

2-(2-methyl-4-chlorophenylamino)-5-(2,6-dimethylphenyl
carbonylamino)pyridine; or

5 2-(2-methyl-4-chlorophenylamino)-5-(2-methylphenyl
aminocarbonyl)pyridine.

16. A pharmaceutical composition comprising a compound
10 of Claim 1 and a pharmaceutically acceptable carrier.

17. A method for prophylaxis or treatment of
inflammation comprising administering an effective
15 amount of a compound of Claim 1.

18. A method for prophylaxis or treatment of
inflammation comprising administering an effective
20 amount of a composition of Claim 16.

19. A method of treating cancer which comprises
administering an effective amount of a compound of Claim
25 1.

20. A method of treating cancer which comprises
administering an effective amount of a composition of
Claim 16.
30

21. A method for prophylaxis or treatment of rheumatoid
arthritis, Pagets disease, osteoporosis, multiple
myeloma, uveitis, acute or chronic myelogenous
35 leukemia, pancreatic β cell destruction, osteoarthritis,
rheumatoid spondylitis, gouty arthritis, inflammatory
bowel disease, adult respiratory distress syndrome

(ARDS), psoriasis, Crohn's disease, allergic rhinitis, ulcerative colitis, anaphylaxis, contact dermatitis, asthma, muscle degeneration, cachexia, Reiter's syndrome, type I diabetes, type II diabetes, bone
5 resorption diseases, graft vs. host reaction, Alzheimer's disease, stroke, myocardial infarction, ischemia reperfusion injury, atherosclerosis, brain trauma, multiple sclerosis, cerebral malaria, sepsis, septic shock, toxic shock syndrome, fever, myalgias due
10 to HIV-1, HIV-2, HIV-3, cytomegalovirus (CMV), influenza, adenovirus, the herpes viruses or herpes zoster infection in a mammal comprising administering an effective amount of a compound of Claim 1.

15

22. A method for prophylaxis or treatment of rheumatoid arthritis, Pagets disease, osteoporosis, multiple myeloma, uveitis, acute or chronic myelogenous leukemia, pancreatic β cell destruction, osteoarthritis,
20 rheumatoid spondylitis, gouty arthritis, inflammatory bowel disease, adult respiratory distress syndrome (ARDS), psoriasis, Crohn's disease, allergic rhinitis, ulcerative colitis, anaphylaxis, contact dermatitis, asthma, muscle degeneration, cachexia, Reiter's
25 syndrome, type I diabetes, type II diabetes, cancer, bone resorption diseases, graft vs. host reaction, Alzheimer's disease, stroke, myocardial infarction, ischemia reperfusion injury, atherosclerosis, brain trauma, multiple sclerosis, cerebral malaria, sepsis,
30 septic shock, toxic shock syndrome, fever, myalgias due to HIV-1, HIV-2, HIV-3, cytomegalovirus (CMV), influenza, adenovirus, the herpes viruses or herpes zoster infection in a mammal comprising administering an effective amount of a compound of Claim 16.

35

23. A method for prophylaxis or treatment of pain comprising administering an effective amount of a compound of Claim 1.

5

24. A method for prophylaxis or treatment of pain comprising administering an effective amount of a composition of Claim 16.

INTERNATIONAL SEARCH REPORT

Inter: nal Application No
PCT/US 98/23510

A. CLASSIFICATION OF SUBJECT MATTER

IPC 6 C07D213/73 C07D213/75 C07D213/82 A61K31/44 C07D413/12
 C07D401/12 C07D409/12 C07D417/12 C07D471/04 C07D405/12

According to International Patent Classification (IPC) or to both national classification and IPC

B. FIELDS SEARCHED

Minimum documentation searched (classification system followed by classification symbols)

IPC 6 C07D A61K

Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched

Electronic data base consulted during the international search (name of data base and, where practical, search terms used)

C. DOCUMENTS CONSIDERED TO BE RELEVANT

| Category * | Citation of document, with indication, where appropriate, of the relevant passages | Relevant to claim No. |
|------------|--|-----------------------|
| X | DE 38 04 346 A (BOEHRINGER MANNHEIM GMBH) 24 August 1989 see the whole document --- | 1, 16 |
| X | EP 0 480 258 A (BAYER AG) 15 April 1992 see claims; example 85 --- | 1, 16 |
| X | DE 18 10 162 A (MERCK & CO., INC.) 14 August 1969 see claims; examples --- | 1, 16 |
| X | EP 0 799 825 A (NIHON NOHYAKU CO LTD) 8 October 1997 see claims; examples --- | 1 |
| X | US 5 380 734 A (HSU KUO-HOM L ET AL) 10 January 1995 see column 6, line 20 --- | 1 |
| -/-- | | |



Further documents are listed in the continuation of box C.



Patent family members are listed in annex.

* Special categories of cited documents :

- "A" document defining the general state of the art which is not considered to be of particular relevance
- "E" earlier document but published on or after the international filing date
- "L" document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other special reason (as specified)
- "O" document referring to an oral disclosure, use, exhibition or other means
- "P" document published prior to the international filing date but later than the priority date claimed

- "T" later document published after the international filing date or priority date and not in conflict with the application but cited to understand the principle or theory underlying the invention
- "X" document of particular relevance; the claimed invention cannot be considered novel or cannot be considered to involve an inventive step when the document is taken alone
- "Y" document of particular relevance; the claimed invention cannot be considered to involve an inventive step when the document is combined with one or more other such documents, such combination being obvious to a person skilled in the art.
- "&" document member of the same patent family

Date of the actual completion of the international search

19 February 1999

Date of mailing of the international search report

15/03/1999

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INTERNATIONAL SEARCH REPORT

Inte: onal Application No
PCT/US 98/23510

C.(Continuation) DOCUMENTS CONSIDERED TO BE RELEVANT

| Category ° | Citation of document, with indication, where appropriate, of the relevant passages | Relevant to claim No. |
|------------|--|-----------------------|
| A | <p>EP 0 424 848 A (HOECHST ROUSSEL PHARMA) 2 May 1991 see the whole document -----</p> | 1, 16-24 |

INTERNATIONAL SEARCH REPORT

International application No.

PCT/US 98/ 23510

Box I Observations where certain claims were found unsearchable (Continuation of item 1 of first sheet)

This International Search Report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:

1. ☒ Claims Nos.: 17-24
because they relate to subject matter not required to be searched by this Authority, namely:
Remark: Although claims 17-24
are directed to a method of treatment of the human/animal
body, the search has been carried out and based on the alleged
effects of the compound/composition.
2. ☐ Claims Nos.:
because they relate to parts of the International Application that do not comply with the prescribed requirements to such
an extent that no meaningful International Search can be carried out, specifically:
3. ☐ Claims Nos.:
because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).

Box II Observations where unity of invention is lacking (Continuation of item 2 of first sheet)

This International Searching Authority found multiple inventions in this international application, as follows:

1. ☐ As all required additional search fees were timely paid by the applicant, this International Search Report covers all
searchable claims.
2. ☐ As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment
of any additional fee.
3. ☐ As only some of the required additional search fees were timely paid by the applicant, this International Search Report
covers only those claims for which fees were paid, specifically claims Nos.:
4. ☐ No required additional search fees were timely paid by the applicant. Consequently, this International Search Report is
restricted to the invention first mentioned in the claims; it is covered by claims Nos.:

Remark on Protest

- ☐ The additional search fees were accompanied by the applicant's protest.
- ☐ No protest accompanied the payment of additional search fees.

INTERNATIONAL SEARCH REPORT

Information on patent family members

International Application No

PCT/US 98/23510

| Patent document cited in search report | | Publication date | Patent family member(s) | Publication date |
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